



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 18, 2024 – 12:42 AM EST

PDB ID : 4I9I  
Title : Crystal structure of tankyrase 1 with compound 4  
Authors : Huang, X.  
Deposited on : 2012-12-05  
Resolution : 2.40 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

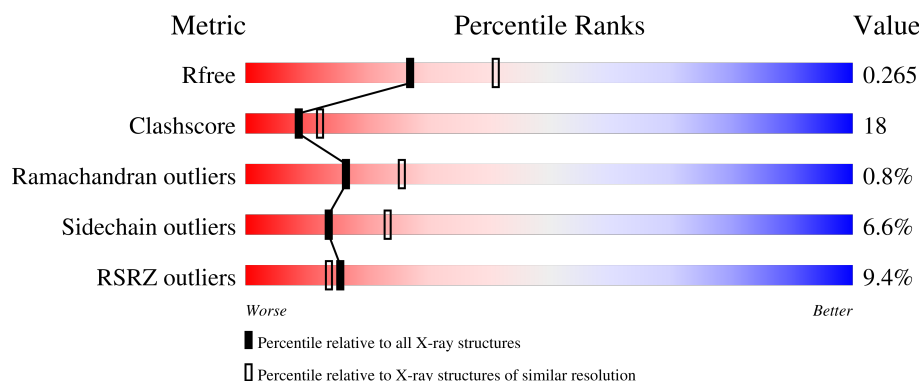
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	164625	4642 (2.40-2.40)
Clashscore	180529	5218 (2.40-2.40)
Ramachandran outliers	177936	5158 (2.40-2.40)
Sidechain outliers	177891	5159 (2.40-2.40)
RSRZ outliers	164620	4642 (2.40-2.40)

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  $\geq 3$ , 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	217	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div>..</div> </div> </div>
1	B	217	<div> <div>9%</div> <div> <div></div> <div>68%</div> <div>25%</div> <div>..</div> </div> </div>
1	C	217	<div> <div>10%</div> <div> <div></div> <div>61%</div> <div>29%</div> <div>5% ..</div> </div> </div>
1	D	217	<div> <div>11%</div> <div> <div></div> <div>61%</div> <div>30%</div> <div>.. 5%</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7021 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 Tankyrase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1698	1066	313	307	12			
1	B	211	Total	C	N	O	S	0	0	0
			1687	1060	310	306	11			
1	C	209	Total	C	N	O	S	0	0	0
			1675	1054	307	303	11			
1	D	206	Total	C	N	O	S	0	0	0
			1651	1038	304	298	11			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1315	HIS	-	expression tag	UNP O95271
A	1316	HIS	-	expression tag	UNP O95271
A	1317	HIS	-	expression tag	UNP O95271
A	1318	HIS	-	expression tag	UNP O95271
A	1319	HIS	-	expression tag	UNP O95271
A	1320	HIS	-	expression tag	UNP O95271
B	1315	HIS	-	expression tag	UNP O95271
B	1316	HIS	-	expression tag	UNP O95271
B	1317	HIS	-	expression tag	UNP O95271
B	1318	HIS	-	expression tag	UNP O95271
B	1319	HIS	-	expression tag	UNP O95271
B	1320	HIS	-	expression tag	UNP O95271
C	1315	HIS	-	expression tag	UNP O95271
C	1316	HIS	-	expression tag	UNP O95271
C	1317	HIS	-	expression tag	UNP O95271
C	1318	HIS	-	expression tag	UNP O95271
C	1319	HIS	-	expression tag	UNP O95271
C	1320	HIS	-	expression tag	UNP O95271
D	1315	HIS	-	expression tag	UNP O95271
D	1316	HIS	-	expression tag	UNP O95271
D	1317	HIS	-	expression tag	UNP O95271

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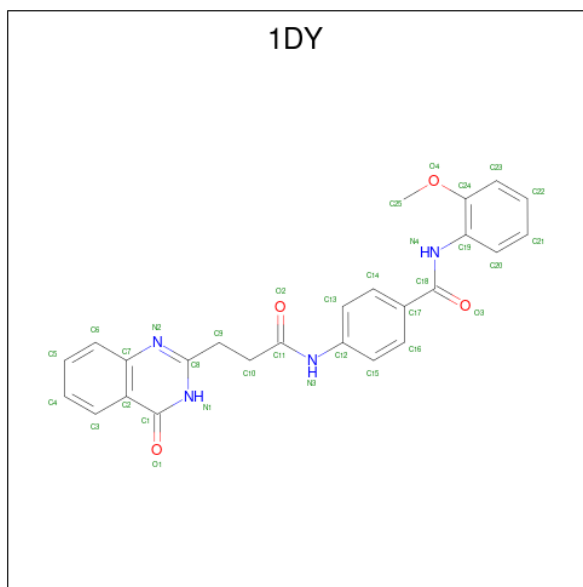
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Chain	Residue	Modelled	Actual	Comment	Reference
D	1318	HIS	-	expression tag	UNP O95271
D	1319	HIS	-	expression tag	UNP O95271
D	1320	HIS	-	expression tag	UNP O95271

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is N-(2-methoxyphenyl)-4-{[3-(4-oxo-3,4-dihydroquinazolin-2-yl)propanoyl]amino}benzamide (three-letter code: 1DY) (formula: C<sub>25</sub>H<sub>22</sub>N<sub>4</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 33 25 4 4	0	0
3	B	1	Total C N O 33 25 4 4	0	0
3	C	1	Total C N O 33 25 4 4	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	D	1	Total	C	N	O	0	0
			33	25	4	4		

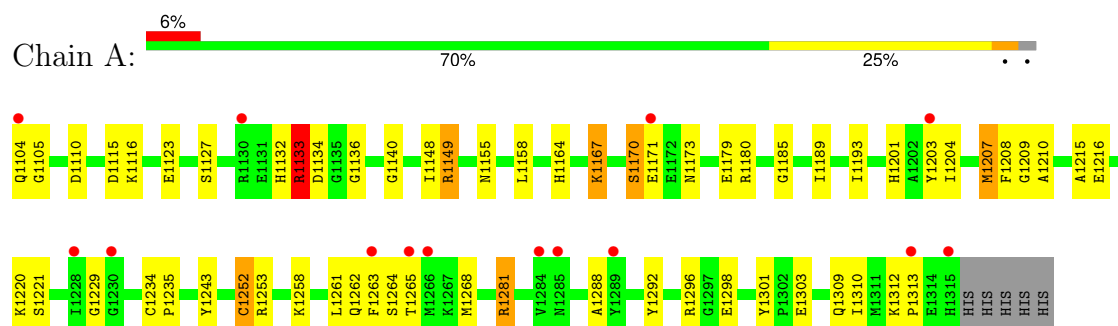
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total	O	0	0
			40	40		
4	B	36	Total	O	0	0
			36	36		
4	C	49	Total	O	0	0
			49	49		
4	D	49	Total	O	0	0
			49	49		

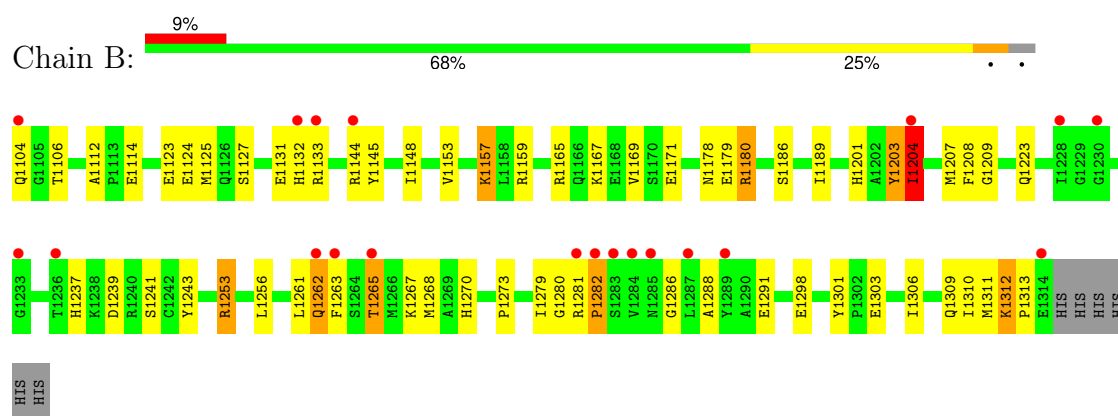
### 3 Residue-property plots [i](#)

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.

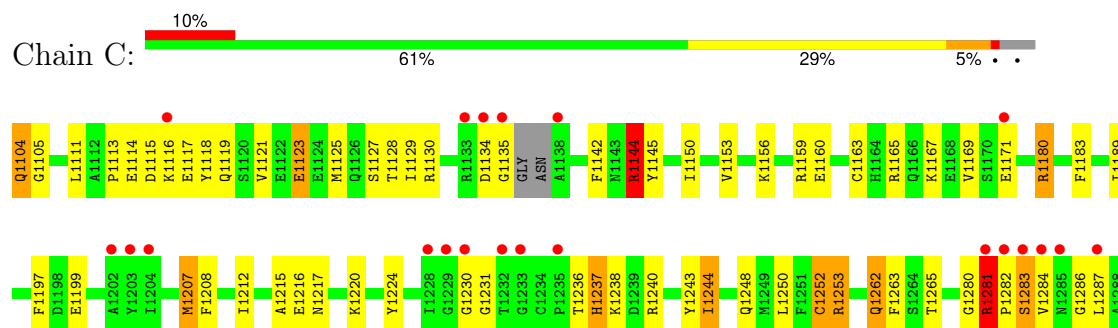
#### • Molecule 1: Tankyrase-1

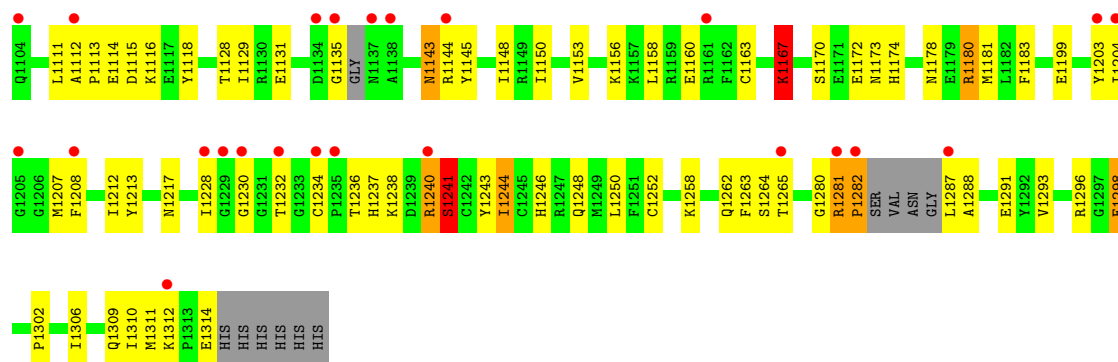


#### • Molecule 1: Tankyrase-1



#### • Molecule 1: Tankyrase-1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	158.62Å 77.12Å 84.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.41	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.40) 98.0 (50.00-2.41)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.99 (at 2.42Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.250 , 0.286 0.236 , 0.265	Depositor DCC
$R_{free}$ test set	2004 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.5	Xtriage
Anisotropy	0.770	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 30.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7021	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.46 % 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 1.7579e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 1DY, ZN

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	0.53	0/1741	0.78	0/2340
1	B	0.53	0/1729	0.85	2/2325 (0.1%)
1	C	0.57	0/1716	0.87	6/2306 (0.3%)
1	D	0.53	0/1691	0.88	6/2272 (0.3%)
All	All	0.54	0/6877	0.84	14/9243 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	1167	LYS	CD-CE-NZ	7.84	129.73	111.70
1	D	1262	GLN	CB-CG-CD	-7.52	92.05	111.60
1	C	1144	ARG	NE-CZ-NH1	6.83	123.71	120.30
1	C	1262	GLN	CB-CG-CD	-6.75	94.04	111.60
1	B	1262	GLN	CB-CG-CD	-6.71	94.16	111.60
1	D	1291	GLU	CB-CG-CD	-6.55	96.52	114.20
1	C	1159	ARG	NE-CZ-NH1	-6.14	117.23	120.30
1	C	1253	ARG	NE-CZ-NH1	-6.08	117.26	120.30
1	B	1253	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	D	1240	ARG	N-CA-C	-5.87	95.16	111.00
1	C	1281	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	D	1172	GLU	CB-CG-CD	-5.49	99.38	114.20
1	D	1135	GLY	N-CA-C	-5.10	100.35	113.10
1	C	1291	GLU	CB-CG-CD	-5.05	100.55	114.20

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	1698	0	1623	50	0
1	B	1687	0	1616	68	0
1	C	1675	0	1606	79	0
1	D	1651	0	1571	55	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	33	0	22	3	0
3	B	33	0	22	5	0
3	C	33	0	22	2	0
3	D	33	0	22	1	0
4	A	40	0	0	4	0
4	B	36	0	0	5	0
4	C	49	0	0	3	0
4	D	49	0	0	5	0
All	All	7021	0	6504	245	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 18.

All (245) 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:C:1281:ARG:NH1	1:C:1281:ARG:HG2	1.58	1.03
1:C:1281:ARG:HG2	1:C:1281:ARG:HH11	0.86	1.01
1:B:1157:LYS:H	1:B:1157:LYS:HZ2	1.03	0.97
1:B:1203:TYR:O	1:B:1204:ILE:HG23	1.65	0.96
1:C:1144:ARG:HG2	1:C:1144:ARG:HH11	1.28	0.96
1:C:1281:ARG:HH11	1:C:1281:ARG:CG	1.79	0.94
1:C:1240:ARG:HG2	1:C:1240:ARG:HH11	1.42	0.85
1:D:1240:ARG:O	1:D:1241:SER:HB2	1.74	0.85
1:B:1180:ARG:HG2	1:B:1180:ARG:HH11	1.48	0.78
1:B:1157:LYS:HB2	1:B:1157:LYS:NZ	2.00	0.75
1:B:1157:LYS:HZ2	1:B:1157:LYS:N	1.82	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1201:HIS:CD2	3:A:1402:1DY:H22	2.21	0.75
1:C:1220:LYS:HE2	1:C:1224:TYR:HE2	1.52	0.73
1:B:1148:ILE:HD11	1:B:1309:GLN:HG3	1.71	0.72
1:A:1179:GLU:OE2	1:A:1253:ARG:NH1	2.22	0.72
1:D:1207:MET:SD	1:D:1287:LEU:HD21	2.30	0.71
1:B:1291:GLU:HG2	3:B:1402:1DY:H4	1.72	0.71
1:B:1167:LYS:O	1:B:1171:GLU:HG3	1.91	0.70
1:A:1167:LYS:O	1:A:1171:GLU:HG3	1.91	0.70
1:A:1262:GLN:NE2	1:A:1268:MET:HE1	2.07	0.69
1:B:1157:LYS:H	1:B:1157:LYS:NZ	1.87	0.69
1:C:1240:ARG:HH11	1:C:1240:ARG:CG	2.04	0.69
1:C:1199:GLU:OE1	1:C:1199:GLU:N	2.23	0.68
1:A:1148:ILE:HD11	1:A:1309:GLN:HG3	1.76	0.68
1:B:1180:ARG:HH11	1:B:1180:ARG:CG	2.06	0.67
1:A:1243:TYR:CE2	1:A:1312:LYS:HE2	2.30	0.67
1:C:1265:THR:O	1:C:1265:THR:HG22	1.95	0.67
1:C:1220:LYS:HE2	1:C:1224:TYR:CE2	2.29	0.66
1:C:1207:MET:HG3	1:C:1265:THR:OG1	1.95	0.66
1:B:1203:TYR:CB	4:B:1507:HOH:O	2.44	0.66
1:C:1314:GLU:HB2	4:C:1544:HOH:O	1.96	0.65
1:B:1180:ARG:HD3	1:C:1281:ARG:NH2	2.11	0.65
1:B:1123:GLU:OE2	1:C:1130:ARG:HA	1.96	0.64
1:D:1163:CYS:SG	4:D:1519:HOH:O	2.54	0.64
1:D:1265:THR:HG22	1:D:1265:THR:O	1.96	0.64
1:C:1189:ILE:HD12	1:C:1250:LEU:HG	1.79	0.64
1:B:1157:LYS:N	1:B:1157:LYS:HD3	2.13	0.64
1:B:1178:ASN:HD21	1:B:1180:ARG:HH12	1.44	0.63
1:B:1208:PHE:CE2	1:B:1280:GLY:HA3	2.34	0.63
1:A:1209:GLY:HA3	1:A:1268:MET:O	1.98	0.63
1:C:1144:ARG:HG2	1:C:1144:ARG:NH1	2.03	0.62
1:C:1244:ILE:HD11	1:D:1236:THR:HG21	1.80	0.62
1:A:1262:GLN:HE21	1:A:1268:MET:HE1	1.65	0.62
1:C:1281:ARG:NH1	1:C:1282:PRO:HG2	2.14	0.62
1:D:1113:PRO:HA	1:D:1118:TYR:CD2	2.34	0.62
1:D:1240:ARG:O	1:D:1241:SER:CB	2.46	0.62
1:B:1203:TYR:C	1:B:1204:ILE:HG23	2.20	0.61
1:D:1128:THR:HB	1:D:1217:ASN:HA	1.82	0.61
1:A:1104:GLN:HG2	1:A:1105:GLY:N	2.14	0.61
1:B:1157:LYS:NZ	1:B:1157:LYS:CB	2.63	0.61
1:C:1119:GLN:O	1:C:1123:GLU:HB2	2.00	0.61
1:B:1243:TYR:CE2	1:B:1312:LYS:HG2	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1296:ARG:HB3	1:D:1298:GLU:OE2	2.01	0.61
1:C:1252:CYS:HB3	1:C:1301:TYR:O	2.01	0.61
1:D:1232:THR:HG23	1:D:1232:THR:O	2.00	0.60
1:A:1261:LEU:HD23	1:A:1281:ARG:HE	1.65	0.60
1:B:1208:PHE:CZ	1:B:1280:GLY:HA3	2.36	0.60
1:C:1180:ARG:HH11	1:C:1180:ARG:HG3	1.66	0.60
1:B:1157:LYS:HB2	1:B:1157:LYS:HZ3	1.64	0.60
1:D:1144:ARG:HG2	1:D:1145:TYR:N	2.16	0.60
1:B:1207:MET:HG2	1:B:1265:THR:OG1	2.01	0.59
1:A:1148:ILE:O	1:A:1149:ARG:HB3	2.00	0.59
1:A:1207:MET:HG3	1:A:1265:THR:OG1	2.03	0.59
1:A:1104:GLN:HG2	1:A:1105:GLY:H	1.65	0.59
1:D:1143:ASN:OD1	1:D:1143:ASN:O	2.20	0.59
1:C:1263:PHE:N	1:C:1263:PHE:CD2	2.70	0.59
1:D:1148:ILE:HD11	1:D:1309:GLN:HG3	1.85	0.58
1:B:1180:ARG:HG3	1:B:1256:LEU:HD12	1.87	0.57
1:C:1281:ARG:NH1	1:C:1281:ARG:CG	2.43	0.57
1:B:1265:THR:HG22	1:B:1265:THR:O	2.04	0.57
1:A:1104:GLN:CG	1:A:1105:GLY:N	2.68	0.57
1:A:1203:TYR:O	1:A:1210:ALA:HA	2.05	0.57
1:A:1133:ARG:CZ	1:A:1288:ALA:HB2	2.35	0.57
1:B:1144:ARG:HG2	1:B:1145:TYR:N	2.19	0.57
1:B:1179:GLU:OE2	1:B:1253:ARG:HD3	2.05	0.57
1:B:1180:ARG:CG	1:B:1180:ARG:NH1	2.65	0.57
1:B:1203:TYR:O	1:B:1204:ILE:CG2	2.48	0.56
1:B:1112:ALA:HB1	1:B:1114:GLU:OE2	2.06	0.56
1:B:1133:ARG:NE	1:B:1288:ALA:HB2	2.21	0.56
1:B:1298:GLU:CD	1:B:1298:GLU:H	2.09	0.56
1:C:1128:THR:HB	1:C:1217:ASN:HA	1.88	0.56
1:D:1212:ILE:HG23	3:D:1402:1DY:H13	1.87	0.56
1:B:1201:HIS:CD2	3:B:1402:1DY:H22	2.42	0.55
1:B:1253:ARG:HG3	1:B:1303:GLU:HG3	1.87	0.55
1:C:1197:PHE:HB3	1:C:1212:ILE:HD13	1.89	0.54
1:B:1279:ILE:HG22	1:B:1281:ARG:HG2	1.90	0.54
1:C:1240:ARG:CG	1:C:1240:ARG:NH1	2.68	0.54
1:A:1132:HIS:C	1:A:1134:ASP:H	2.10	0.54
1:C:1153:VAL:HB	1:C:1302:PRO:O	2.08	0.54
1:C:1129:ILE:HD11	1:C:1145:TYR:CE2	2.43	0.54
1:A:1203:TYR:CB	4:A:1503:HOH:O	2.57	0.53
1:D:1129:ILE:HD11	1:D:1145:TYR:CE2	2.43	0.53
1:B:1157:LYS:HZ2	1:B:1157:LYS:HB2	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1263:PHE:N	1:A:1263:PHE:CD2	2.77	0.52
1:C:1243:TYR:CE2	1:C:1312:LYS:HG3	2.44	0.52
1:C:1180:ARG:NH2	1:C:1216:GLU:OE1	2.42	0.52
1:A:1173:ASN:ND2	1:A:1258:LYS:HB2	2.24	0.52
1:C:1134:ASP:HB2	1:C:1289:TYR:HE1	1.75	0.52
1:D:1207:MET:CE	1:D:1265:THR:OG1	2.58	0.52
1:D:1243:TYR:CD2	1:D:1312:LYS:HD3	2.45	0.52
1:C:1281:ARG:NH2	4:C:1549:HOH:O	2.43	0.51
1:D:1180:ARG:NE	4:D:1532:HOH:O	2.32	0.51
1:A:1180:ARG:NH2	1:A:1216:GLU:OE1	2.43	0.51
1:B:1132:HIS:CD2	1:B:1223:GLN:HE21	2.28	0.51
1:D:1232:THR:O	1:D:1238:LYS:HA	2.10	0.51
1:D:1234:CYS:SG	1:D:1237:HIS:N	2.83	0.51
1:A:1123:GLU:O	1:A:1127:SER:OG	2.28	0.51
1:D:1228:ILE:C	1:D:1230:GLY:H	2.14	0.51
1:A:1164:HIS:CD2	4:A:1539:HOH:O	2.63	0.51
1:C:1253:ARG:HG3	1:C:1303:GLU:HG3	1.92	0.51
1:C:1115:ASP:OD1	1:C:1116:LYS:N	2.44	0.51
1:C:1237:HIS:CD2	1:D:1237:HIS:CD2	2.99	0.51
1:D:1199:GLU:H	1:D:1199:GLU:CD	2.13	0.51
1:A:1207:MET:HB3	1:A:1208:PHE:CD2	2.45	0.50
1:D:1144:ARG:HD3	1:D:1311:MET:SD	2.51	0.50
1:A:1115:ASP:OD1	1:A:1116:LYS:N	2.44	0.50
1:A:1265:THR:O	1:A:1265:THR:HG22	2.11	0.50
1:A:1132:HIS:O	1:A:1134:ASP:N	2.45	0.50
1:A:1155:ASN:HB3	1:A:1158:LEU:HB2	1.93	0.50
1:C:1281:ARG:HH12	1:C:1282:PRO:HG2	1.76	0.50
1:B:1157:LYS:HZ2	1:B:1157:LYS:CB	2.24	0.50
1:C:1282:PRO:HA	1:C:1287:LEU:HB2	1.93	0.50
1:C:1113:PRO:HA	1:C:1118:TYR:CG	2.46	0.50
1:C:1113:PRO:HA	1:C:1118:TYR:CD2	2.47	0.49
1:C:1237:HIS:O	1:C:1238:LYS:C	2.49	0.49
1:D:1112:ALA:O	1:D:1115:ASP:HB2	2.12	0.49
1:C:1283:SER:HB3	1:C:1286:GLY:HA3	1.94	0.49
1:B:1291:GLU:HG2	3:B:1402:1DY:C5	2.43	0.49
1:C:1142:PHE:HA	1:C:1313:PRO:HG3	1.93	0.49
1:C:1292:TYR:CD2	1:C:1292:TYR:N	2.81	0.49
1:D:1113:PRO:HA	1:D:1118:TYR:CG	2.48	0.49
1:D:1234:CYS:SG	1:D:1238:LYS:N	2.86	0.49
1:C:1114:GLU:CD	1:C:1114:GLU:H	2.16	0.48
1:C:1281:ARG:CG	1:C:1282:PRO:HD2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1129:ILE:HD11	1:D:1145:TYR:CD2	2.49	0.48
1:D:1167:LYS:O	1:D:1167:LYS:HG3	2.13	0.48
1:B:1144:ARG:HG2	1:B:1145:TYR:H	1.79	0.48
1:B:1273:PRO:HD2	4:B:1531:HOH:O	2.13	0.48
1:C:1115:ASP:OD1	1:C:1115:ASP:C	2.52	0.48
1:C:1144:ARG:NH1	1:C:1144:ARG:CG	2.76	0.48
1:A:1253:ARG:HG3	1:A:1303:GLU:HG3	1.96	0.48
1:A:1132:HIS:C	1:A:1134:ASP:N	2.67	0.48
1:A:1167:LYS:HA	1:A:1170:SER:HB2	1.95	0.48
1:C:1129:ILE:HD11	1:C:1145:TYR:CD2	2.49	0.48
1:D:1213:TYR:CD1	1:D:1293:VAL:HG22	2.50	0.47
1:B:1133:ARG:NH1	1:B:1286:GLY:O	2.48	0.47
1:B:1133:ARG:CZ	1:B:1288:ALA:HB2	2.44	0.47
1:D:1180:ARG:HG3	1:D:1180:ARG:HH11	1.80	0.47
1:A:1268:MET:HE3	1:A:1268:MET:HB2	1.74	0.47
1:C:1156:LYS:O	1:C:1160:GLU:HG3	2.14	0.47
1:C:1283:SER:CB	1:C:1286:GLY:HA3	2.45	0.47
1:D:1180:ARG:HG3	4:D:1544:HOH:O	2.15	0.47
1:B:1270:HIS:HD2	4:B:1516:HOH:O	1.97	0.47
1:C:1167:LYS:HE3	1:C:1167:LYS:HB3	1.77	0.47
1:B:1243:TYR:CD1	1:B:1310:ILE:HG13	2.49	0.46
1:B:1239:ASP:OD1	1:B:1241:SER:N	2.31	0.46
1:C:1134:ASP:CG	1:C:1135:GLY:H	2.19	0.46
1:C:1281:ARG:HG2	1:C:1282:PRO:HD2	1.98	0.46
1:D:1112:ALA:HB1	1:D:1114:GLU:OE2	2.16	0.46
1:A:1215:ALA:HB2	3:A:1402:1DY:H3	1.98	0.46
1:C:1215:ALA:HB2	3:C:1402:1DY:H3	1.98	0.46
1:D:1263:PHE:N	1:D:1263:PHE:CD2	2.83	0.46
1:B:1106:THR:HG21	1:B:1153:VAL:HG13	1.96	0.46
1:B:1267:LYS:HB2	4:B:1533:HOH:O	2.15	0.45
1:B:1301:TYR:HD2	1:B:1303:GLU:HG2	1.82	0.45
1:C:1207:MET:HB3	1:C:1208:PHE:CE2	2.51	0.45
1:C:1236:THR:HB	1:C:1237:HIS:CE1	2.51	0.45
1:D:1111:LEU:HD12	1:D:1150:ILE:HG22	1.99	0.45
1:B:1263:PHE:N	1:B:1263:PHE:CD2	2.84	0.45
1:C:1248:GLN:HA	1:C:1306:ILE:O	2.15	0.45
1:D:1207:MET:HE2	1:D:1265:THR:OG1	2.16	0.45
1:C:1230:GLY:O	1:C:1231:GLY:C	2.55	0.45
1:A:1207:MET:HB3	1:A:1208:PHE:CE2	2.52	0.45
1:D:1144:ARG:HG2	1:D:1145:TYR:H	1.82	0.45
1:D:1183:PHE:HA	1:D:1250:LEU:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1127:SER:HB2	1:C:1127:SER:O	2.16	0.45
1:B:1243:TYR:HD1	1:B:1310:ILE:O	2.00	0.45
1:B:1262:GLN:HE21	1:B:1262:GLN:HB3	1.29	0.45
1:C:1114:GLU:O	1:C:1119:GLN:NE2	2.50	0.45
1:D:1111:LEU:HD12	1:D:1150:ILE:CG2	2.47	0.45
1:D:1153:VAL:HB	1:D:1302:PRO:O	2.17	0.45
1:D:1248:GLN:HA	1:D:1306:ILE:O	2.17	0.44
1:D:1310:ILE:O	1:D:1310:ILE:HG13	2.17	0.44
1:A:1243:TYR:CD1	1:A:1310:ILE:HG13	2.53	0.44
1:C:1281:ARG:NH1	1:C:1282:PRO:CG	2.78	0.44
1:B:1239:ASP:OD1	1:B:1239:ASP:C	2.55	0.44
1:C:1281:ARG:HH11	1:C:1282:PRO:HD2	1.82	0.44
1:B:1186:SER:O	1:B:1189:ILE:HG12	2.17	0.44
1:C:1262:GLN:O	1:C:1280:GLY:HA2	2.18	0.44
1:C:1134:ASP:CG	1:C:1135:GLY:N	2.72	0.43
1:C:1104:GLN:HG2	1:C:1105:GLY:N	2.29	0.43
1:D:1113:PRO:HB3	1:D:1118:TYR:CZ	2.53	0.43
1:C:1165:ARG:O	1:C:1169:VAL:HG23	2.19	0.43
1:D:1208:PHE:CE2	1:D:1280:GLY:HA3	2.54	0.43
1:D:1281:ARG:NH1	4:D:1535:HOH:O	2.52	0.43
1:C:1134:ASP:HB2	1:C:1289:TYR:CE1	2.53	0.43
1:C:1180:ARG:HH11	1:C:1180:ARG:CG	2.31	0.43
1:A:1136:GLY:O	1:A:1140:GLY:N	2.52	0.43
1:A:1185:GLY:N	1:A:1221:SER:HB3	2.34	0.43
1:D:1181:MET:HA	1:D:1252:CYS:O	2.18	0.43
1:A:1252:CYS:HB3	1:A:1301:TYR:O	2.19	0.42
1:B:1159:ARG:NH1	4:B:1510:HOH:O	2.52	0.42
1:D:1244:ILE:O	1:D:1246:HIS:HD2	2.02	0.42
1:A:1189:ILE:O	1:A:1193:ILE:HG12	2.18	0.42
1:C:1117:GLU:O	1:C:1121:VAL:HG23	2.19	0.42
1:D:1156:LYS:O	1:D:1160:GLU:HG3	2.19	0.42
1:D:1265:THR:O	1:D:1265:THR:CG2	2.67	0.42
1:A:1201:HIS:CG	3:A:1402:1DY:H22	2.53	0.42
1:A:1220:LYS:HA	1:A:1220:LYS:HD2	1.85	0.42
1:A:1312:LYS:HA	1:A:1313:PRO:HD3	1.71	0.42
1:B:1165:ARG:O	1:B:1169:VAL:HG23	2.19	0.42
1:C:1111:LEU:HD12	1:C:1150:ILE:CG2	2.49	0.42
1:A:1180:ARG:NH1	1:A:1292:TYR:OH	2.51	0.42
1:A:1229:GLY:N	4:A:1515:HOH:O	2.53	0.42
1:B:1291:GLU:CG	3:B:1402:1DY:H4	2.46	0.42
1:D:1174:HIS:CD2	1:D:1174:HIS:N	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1243:TYR:HD1	1:B:1310:ILE:HG13	1.84	0.42
1:B:1261:LEU:HG	1:B:1281:ARG:HD2	2.01	0.42
1:C:1207:MET:O	1:C:1265:THR:HG23	2.20	0.42
1:C:1163:CYS:SG	4:C:1531:HOH:O	2.62	0.42
1:A:1204:ILE:CG2	4:A:1520:HOH:O	2.67	0.42
1:B:1157:LYS:N	1:B:1157:LYS:CD	2.83	0.42
1:B:1209:GLY:HA3	1:B:1268:MET:O	2.20	0.42
1:C:1167:LYS:O	1:C:1171:GLU:HG3	2.20	0.41
1:A:1148:ILE:CD1	1:A:1309:GLN:HG3	2.48	0.41
1:B:1180:ARG:CZ	1:C:1281:ARG:HE	2.33	0.41
1:D:1207:MET:SD	1:D:1287:LEU:CD2	3.06	0.41
1:A:1296:ARG:HB3	1:A:1298:GLU:OE2	2.21	0.41
1:D:1288:ALA:HB2	4:D:1528:HOH:O	2.20	0.41
1:A:1133:ARG:HE	1:A:1133:ARG:HB3	1.46	0.41
1:C:1237:HIS:ND1	1:C:1237:HIS:N	2.69	0.41
1:A:1312:LYS:HD2	1:B:1237:HIS:CD2	2.56	0.41
1:B:1179:GLU:CD	1:B:1253:ARG:HD3	2.41	0.41
1:C:1125:MET:HA	1:C:1183:PHE:CZ	2.56	0.41
1:D:1282:PRO:HA	1:D:1287:LEU:HB2	2.02	0.41
3:B:1402:1DY:H16	3:B:1402:1DY:O3	2.21	0.41
1:C:1253:ARG:HH11	1:C:1253:ARG:HD3	1.70	0.41
1:B:1125:MET:SD	1:B:1306:ILE:HG21	2.61	0.40
1:C:1212:ILE:HG23	3:C:1402:1DY:H13	2.03	0.40
1:A:1234:CYS:HA	1:A:1235:PRO:HD3	1.75	0.40
1:D:1173:ASN:ND2	1:D:1258:LYS:HB2	2.36	0.40
1:B:1133:ARG:HE	1:B:1133:ARG:HB3	1.77	0.40
1:B:1180:ARG:HD3	1:C:1281:ARG:HH21	1.85	0.40
1:D:1207:MET:HE3	1:D:1265:THR:OG1	2.20	0.40
1:B:1312:LYS:HA	1:B:1313:PRO:HD3	1.86	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	210/217 (97%)	204 (97%)	5 (2%)	1 (0%)	25	38
1	B	209/217 (96%)	200 (96%)	6 (3%)	3 (1%)	9	13
1	C	205/217 (94%)	194 (95%)	11 (5%)	0	100	100
1	D	200/217 (92%)	191 (96%)	6 (3%)	3 (2%)	8	12
All	All	824/868 (95%)	789 (96%)	28 (3%)	7 (1%)	16	26

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1203	TYR
1	B	1204	ILE
1	D	1203	TYR
1	D	1204	ILE
1	D	1241	SER
1	A	1133	ARG
1	B	1282	PRO

### 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	177/185 (96%)	168 (95%)	9 (5%)	20	35
1	B	176/185 (95%)	166 (94%)	10 (6%)	17	29
1	C	175/185 (95%)	163 (93%)	12 (7%)	13	22
1	D	171/185 (92%)	156 (91%)	15 (9%)	8	13
All	All	699/740 (94%)	653 (93%)	46 (7%)	14	23

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

Mol	Chain	Res	Type
1	A	1110	ASP
1	A	1133	ARG
1	A	1149	ARG
1	A	1167	LYS

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Mol	Chain	Res	Type
1	A	1170	SER
1	A	1207	MET
1	A	1252	CYS
1	A	1264	SER
1	A	1281	ARG
1	B	1104	GLN
1	B	1124	GLU
1	B	1131	GLU
1	B	1157	LYS
1	B	1180	ARG
1	B	1204	ILE
1	B	1265	THR
1	B	1282	PRO
1	B	1311	MET
1	B	1312	LYS
1	C	1104	GLN
1	C	1123	GLU
1	C	1144	ARG
1	C	1180	ARG
1	C	1207	MET
1	C	1237	HIS
1	C	1244	ILE
1	C	1252	CYS
1	C	1281	ARG
1	C	1283	SER
1	C	1284	VAL
1	C	1314	GLU
1	D	1116	LYS
1	D	1131	GLU
1	D	1143	ASN
1	D	1158	LEU
1	D	1167	LYS
1	D	1170	SER
1	D	1178	ASN
1	D	1180	ARG
1	D	1241	SER
1	D	1244	ILE
1	D	1264	SER
1	D	1281	ARG
1	D	1282	PRO
1	D	1298	GLU
1	D	1314	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1173	ASN
1	A	1201	HIS
1	A	1248	GLN
1	A	1262	GLN
1	B	1104	GLN
1	B	1119	GLN
1	B	1178	ASN
1	B	1201	HIS
1	B	1217	ASN
1	B	1223	GLN
1	B	1262	GLN
1	C	1119	GLN
1	C	1173	ASN
1	C	1223	GLN
1	C	1262	GLN
1	D	1119	GLN
1	D	1166	GLN
1	D	1173	ASN
1	D	1223	GLN
1	D	1246	HIS
1	D	1248	GLN
1	D	1262	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
3	1DY	D	1402	-	34,36,36	2.71	20 (58%)	46,49,49	1.68	12 (26%)
3	1DY	C	1402	-	34,36,36	2.53	21 (61%)	46,49,49	1.69	11 (23%)
3	1DY	A	1402	-	34,36,36	2.52	19 (55%)	46,49,49	2.17	11 (23%)
3	1DY	B	1402	-	34,36,36	2.57	20 (58%)	46,49,49	2.02	14 (30%)

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	1DY	D	1402	-	-	0/19/19/19	0/4/4/4
3	1DY	C	1402	-	-	2/19/19/19	0/4/4/4
3	1DY	A	1402	-	-	3/19/19/19	0/4/4/4
3	1DY	B	1402	-	-	3/19/19/19	0/4/4/4

All (80) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1402	1DY	C14-C17	5.11	1.47	1.39
3	B	1402	1DY	C2-C7	4.65	1.46	1.40
3	A	1402	1DY	C14-C17	4.61	1.46	1.39
3	C	1402	1DY	C14-C13	4.57	1.46	1.38
3	A	1402	1DY	C21-C20	4.55	1.46	1.38
3	D	1402	1DY	C2-C7	4.43	1.46	1.40
3	D	1402	1DY	C14-C17	4.33	1.46	1.39
3	D	1402	1DY	C6-C7	4.31	1.46	1.40
3	B	1402	1DY	C14-C17	4.21	1.45	1.39
3	A	1402	1DY	C14-C13	4.20	1.45	1.38
3	D	1402	1DY	C14-C13	4.18	1.45	1.38
3	B	1402	1DY	C21-C20	4.15	1.46	1.38
3	B	1402	1DY	C14-C13	4.14	1.45	1.38
3	D	1402	1DY	C16-C15	4.12	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1402	1DY	C15-C12	3.94	1.45	1.39
3	D	1402	1DY	C5-C6	3.94	1.45	1.38
3	B	1402	1DY	C15-C12	3.92	1.45	1.39
3	D	1402	1DY	C4-C3	3.83	1.45	1.38
3	A	1402	1DY	C15-C12	3.74	1.45	1.39
3	C	1402	1DY	C16-C15	3.66	1.44	1.38
3	A	1402	1DY	C23-C24	3.50	1.46	1.39
3	D	1402	1DY	C5-C4	3.49	1.45	1.38
3	C	1402	1DY	C5-C6	3.40	1.44	1.38
3	B	1402	1DY	C22-C23	3.36	1.44	1.38
3	C	1402	1DY	C16-C17	3.29	1.44	1.39
3	D	1402	1DY	C16-C17	3.29	1.44	1.39
3	A	1402	1DY	C4-C3	3.28	1.44	1.38
3	C	1402	1DY	C13-C12	3.27	1.44	1.39
3	C	1402	1DY	C23-C24	3.27	1.46	1.39
3	C	1402	1DY	C6-C7	3.23	1.45	1.40
3	A	1402	1DY	C22-C23	3.21	1.44	1.38
3	A	1402	1DY	C2-C7	3.18	1.44	1.40
3	C	1402	1DY	C12-N3	-3.18	1.35	1.41
3	B	1402	1DY	C5-C4	3.17	1.45	1.38
3	D	1402	1DY	C12-N3	-3.17	1.35	1.41
3	B	1402	1DY	C20-C19	3.16	1.44	1.39
3	C	1402	1DY	C21-C20	3.10	1.44	1.38
3	D	1402	1DY	C21-C20	3.08	1.44	1.38
3	B	1402	1DY	C4-C3	3.08	1.44	1.38
3	C	1402	1DY	C5-C4	3.08	1.45	1.38
3	C	1402	1DY	C15-C12	3.07	1.44	1.39
3	B	1402	1DY	C7-N2	-3.05	1.33	1.39
3	B	1402	1DY	C16-C17	3.05	1.44	1.39
3	A	1402	1DY	C1-N1	-3.04	1.33	1.37
3	B	1402	1DY	C16-C15	3.02	1.43	1.38
3	A	1402	1DY	C5-C6	3.00	1.44	1.38
3	D	1402	1DY	C13-C12	2.99	1.44	1.39
3	C	1402	1DY	C20-C19	2.97	1.44	1.39
3	A	1402	1DY	C20-C19	2.96	1.44	1.39
3	A	1402	1DY	C22-C21	2.90	1.44	1.38
3	A	1402	1DY	C6-C7	2.90	1.44	1.40
3	B	1402	1DY	C23-C24	2.90	1.45	1.39
3	B	1402	1DY	C12-N3	-2.81	1.36	1.41
3	D	1402	1DY	C20-C19	2.81	1.44	1.39
3	A	1402	1DY	C13-C12	2.71	1.43	1.39
3	D	1402	1DY	C22-C21	2.70	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1402	1DY	C12-N3	-2.70	1.36	1.41
3	D	1402	1DY	C22-C23	2.66	1.43	1.38
3	C	1402	1DY	C2-C1	-2.60	1.42	1.47
3	B	1402	1DY	C3-C2	2.56	1.43	1.39
3	D	1402	1DY	C23-C24	2.54	1.44	1.39
3	A	1402	1DY	C16-C17	2.52	1.43	1.39
3	A	1402	1DY	C5-C4	2.50	1.43	1.38
3	C	1402	1DY	C2-C7	2.48	1.43	1.40
3	B	1402	1DY	C13-C12	2.46	1.43	1.39
3	B	1402	1DY	C22-C21	2.44	1.43	1.38
3	D	1402	1DY	C3-C2	2.44	1.43	1.39
3	B	1402	1DY	C5-C6	2.42	1.43	1.38
3	A	1402	1DY	C16-C15	2.34	1.42	1.38
3	A	1402	1DY	C7-N2	-2.30	1.35	1.39
3	C	1402	1DY	C9-C8	2.23	1.53	1.49
3	D	1402	1DY	C7-N2	-2.21	1.35	1.39
3	D	1402	1DY	C9-C8	2.21	1.53	1.49
3	C	1402	1DY	C22-C21	2.19	1.43	1.38
3	B	1402	1DY	C1-N1	-2.19	1.34	1.37
3	C	1402	1DY	C19-N4	-2.18	1.37	1.41
3	C	1402	1DY	C4-C3	2.18	1.42	1.38
3	B	1402	1DY	C6-C7	2.18	1.43	1.40
3	C	1402	1DY	C22-C23	2.16	1.42	1.38
3	C	1402	1DY	C19-C24	2.16	1.44	1.40

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1402	1DY	O4-C24-C19	-7.47	105.51	114.81
3	B	1402	1DY	C1-N1-C8	5.92	127.32	123.90
3	A	1402	1DY	C10-C9-C8	-5.39	107.00	112.55
3	D	1402	1DY	C1-N1-C8	4.43	126.46	123.90
3	A	1402	1DY	C16-C17-C14	4.32	124.06	118.57
3	A	1402	1DY	O4-C24-C23	3.98	131.02	124.30
3	B	1402	1DY	C16-C17-C14	3.96	123.61	118.57
3	C	1402	1DY	O1-C1-N1	3.89	125.23	120.62
3	B	1402	1DY	C6-C7-C2	3.71	122.79	119.14
3	A	1402	1DY	C2-C7-N2	-3.64	118.75	122.41
3	A	1402	1DY	C15-C16-C17	-3.59	116.97	120.80
3	C	1402	1DY	C2-C7-N2	-3.52	118.88	122.41
3	B	1402	1DY	O4-C24-C19	-3.47	110.50	114.81
3	D	1402	1DY	C16-C17-C14	3.32	122.80	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1402	1DY	C15-C12-C13	3.31	123.44	119.04
3	C	1402	1DY	N1-C8-N2	-3.29	118.91	123.08
3	D	1402	1DY	C25-O4-C24	3.21	122.22	117.51
3	B	1402	1DY	C15-C12-C13	3.13	123.19	119.04
3	D	1402	1DY	C15-C12-C13	3.09	123.15	119.04
3	B	1402	1DY	N1-C8-N2	-3.05	119.21	123.08
3	D	1402	1DY	C6-C7-C2	3.04	122.13	119.14
3	D	1402	1DY	C2-C7-N2	-2.98	119.42	122.41
3	A	1402	1DY	C15-C12-C13	2.80	122.76	119.04
3	B	1402	1DY	C15-C16-C17	-2.79	117.81	120.80
3	C	1402	1DY	C1-N1-C8	2.73	125.48	123.90
3	A	1402	1DY	C6-C7-C2	2.70	121.80	119.14
3	C	1402	1DY	C3-C2-C7	2.67	122.83	119.78
3	D	1402	1DY	N1-C8-N2	-2.63	119.74	123.08
3	C	1402	1DY	C16-C17-C14	2.59	121.86	118.57
3	B	1402	1DY	C2-C7-N2	-2.53	119.87	122.41
3	C	1402	1DY	C10-C9-C8	-2.52	109.95	112.55
3	B	1402	1DY	C4-C3-C2	-2.47	115.33	119.80
3	B	1402	1DY	C4-C5-C6	2.47	123.28	120.24
3	C	1402	1DY	C12-N3-C11	2.46	131.88	127.52
3	B	1402	1DY	C13-C14-C17	-2.45	118.18	120.80
3	D	1402	1DY	C2-C1-N1	-2.41	112.39	114.93
3	A	1402	1DY	C13-C14-C17	-2.41	118.23	120.80
3	A	1402	1DY	C14-C13-C12	-2.41	117.53	120.30
3	B	1402	1DY	C5-C6-C7	-2.32	114.71	118.97
3	A	1402	1DY	C23-C24-C19	2.28	123.47	120.39
3	D	1402	1DY	C13-C12-N3	-2.18	113.10	120.41
3	D	1402	1DY	O1-C1-C2	2.17	127.14	123.27
3	B	1402	1DY	O4-C24-C23	2.13	127.90	124.30
3	C	1402	1DY	O1-C1-C2	-2.09	119.53	123.27
3	D	1402	1DY	C15-C16-C17	-2.07	118.58	120.80
3	C	1402	1DY	C14-C13-C12	-2.02	117.97	120.30
3	B	1402	1DY	C10-C9-C8	-2.01	110.48	112.55
3	D	1402	1DY	C10-C9-C8	2.01	114.62	112.55

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1402	1DY	N1-C8-C9-C10
3	A	1402	1DY	N2-C8-C9-C10
3	B	1402	1DY	N1-C8-C9-C10

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Mol	Chain	Res	Type	Atoms
3	B	1402	1DY	N2-C8-C9-C10
3	A	1402	1DY	C11-C10-C9-C8
3	B	1402	1DY	C11-C10-C9-C8
3	C	1402	1DY	N1-C8-C9-C10
3	C	1402	1DY	N2-C8-C9-C10

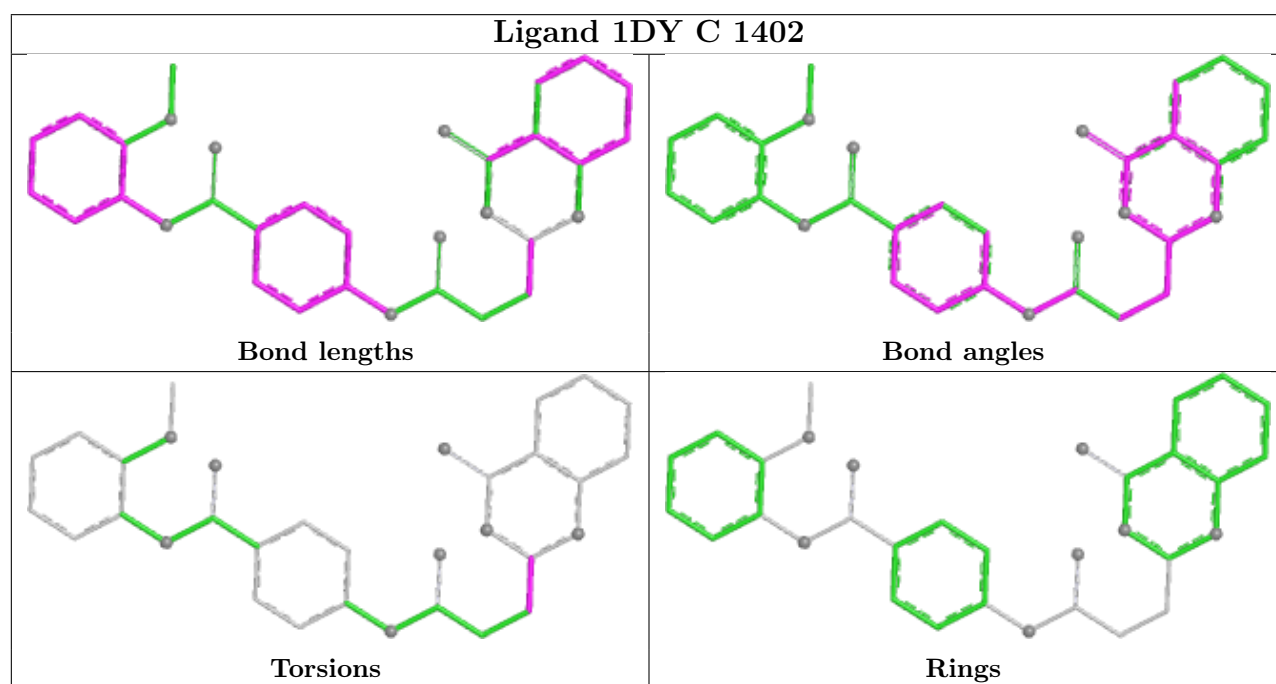
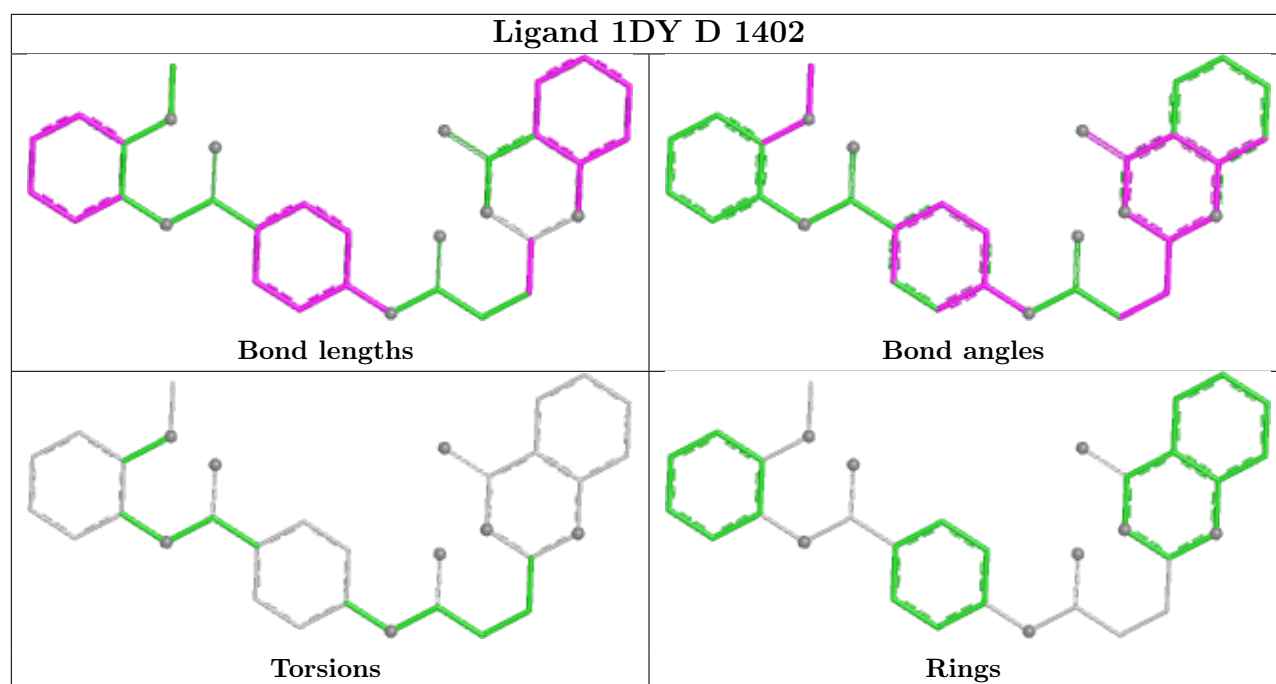
There are no ring outliers.

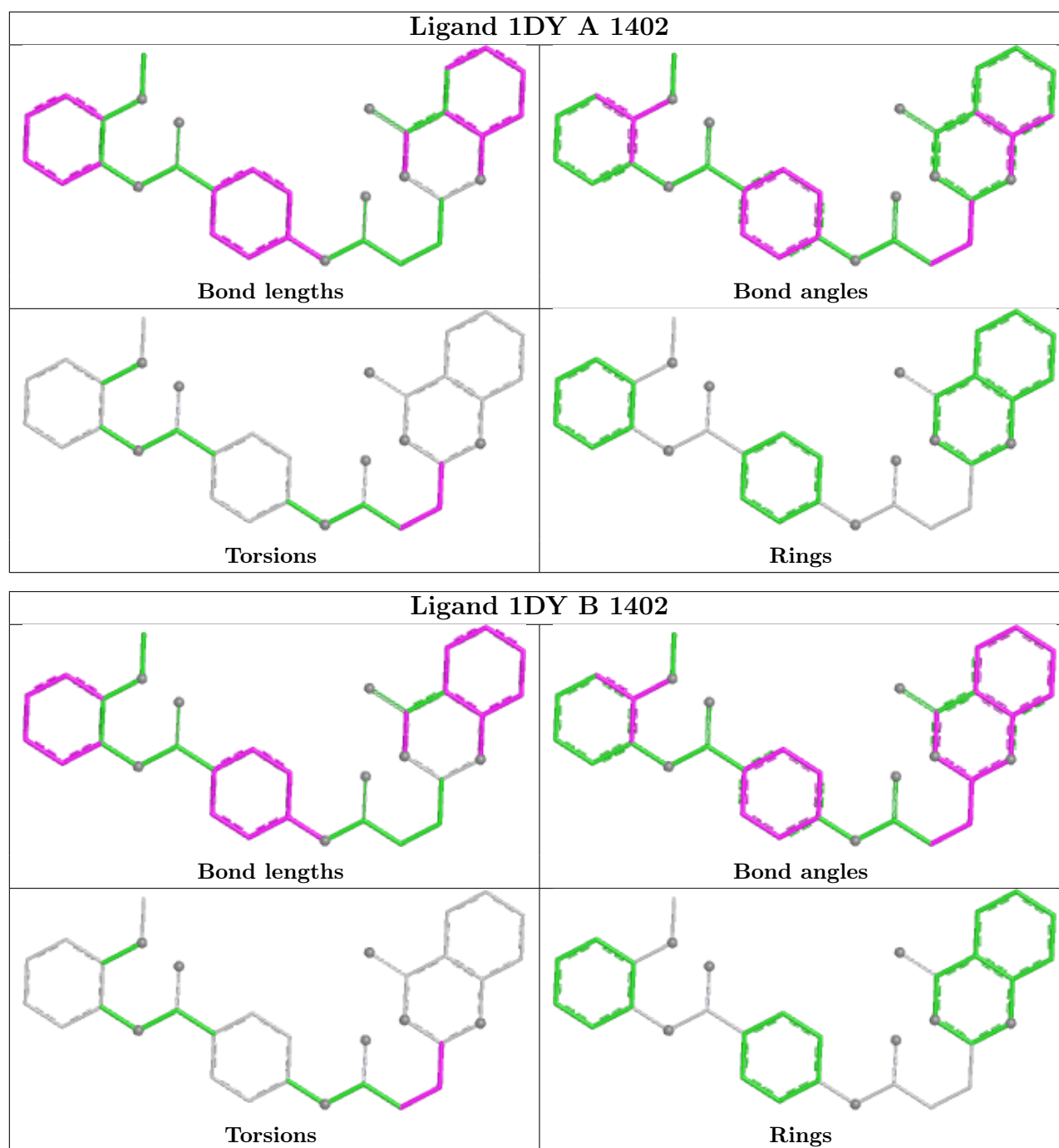
4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1402	1DY	1	0
3	C	1402	1DY	2	0
3	A	1402	1DY	3	0
3	B	1402	1DY	5	0

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







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	212/217 (97%)	0.41	14 (6%) 26 23	21, 37, 63, 82	0
1	B	211/217 (97%)	0.51	20 (9%) 15 13	19, 34, 67, 79	0
1	C	209/217 (96%)	0.62	21 (10%) 14 12	22, 38, 68, 73	0
1	D	206/217 (94%)	0.67	24 (11%) 10 9	22, 38, 67, 75	0
All	All	838/868 (96%)	0.55	79 (9%) 15 14	19, 37, 67, 82	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1203	TYR	5.1
1	B	1285	ASN	4.6
1	C	1230	GLY	4.5
1	D	1282	PRO	4.3
1	C	1285	ASN	4.2
1	B	1282	PRO	4.1
1	D	1232	THR	4.1
1	C	1228	ILE	4.1
1	B	1283	SER	3.7
1	D	1203	TYR	3.7
1	D	1137	ASN	3.7
1	C	1229	GLY	3.6
1	D	1204	ILE	3.6
1	D	1112	ALA	3.6
1	D	1228	ILE	3.5
1	C	1204	ILE	3.5
1	D	1138	ALA	3.5
1	D	1287	LEU	3.4
1	B	1281	ARG	3.4
1	C	1283	SER	3.2
1	C	1281	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	1203	TYR	3.2
1	C	1284	VAL	3.0
1	A	1285	ASN	2.9
1	D	1144	ARG	2.8
1	D	1230	GLY	2.8
1	D	1134	ASP	2.8
1	B	1265	THR	2.8
1	B	1233	GLY	2.8
1	A	1104	GLN	2.8
1	D	1135	GLY	2.8
1	D	1312	LYS	2.8
1	B	1204	ILE	2.7
1	B	1314	GLU	2.7
1	A	1289	TYR	2.7
1	A	1265	THR	2.7
1	D	1205	GLY	2.6
1	B	1284	VAL	2.6
1	B	1228	ILE	2.6
1	A	1266	MET	2.6
1	A	1230	GLY	2.5
1	A	1315	HIS	2.5
1	A	1228	ILE	2.5
1	B	1236	THR	2.5
1	D	1208	PHE	2.5
1	C	1135	GLY	2.4
1	A	1130	ARG	2.4
1	D	1240	ARG	2.4
1	C	1232	THR	2.4
1	C	1233	GLY	2.4
1	C	1134	ASP	2.4
1	A	1284	VAL	2.4
1	C	1287	LEU	2.3
1	B	1144	ARG	2.3
1	D	1234	CYS	2.3
1	B	1262	GLN	2.3
1	B	1132	HIS	2.3
1	D	1265	THR	2.3
1	B	1263	PHE	2.3
1	C	1202	ALA	2.3
1	A	1171	GLU	2.3
1	C	1138	ALA	2.2
1	C	1133	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	1289	TYR	2.2
1	C	1235	PRO	2.2
1	D	1104	GLN	2.2
1	A	1313	PRO	2.1
1	B	1133	ARG	2.1
1	D	1235	PRO	2.1
1	C	1171	GLU	2.1
1	D	1161	ARG	2.1
1	C	1282	PRO	2.1
1	C	1116	LYS	2.1
1	D	1281	ARG	2.1
1	B	1230	GLY	2.1
1	A	1263	PHE	2.1
1	B	1287	LEU	2.0
1	D	1229	GLY	2.0
1	B	1104	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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.

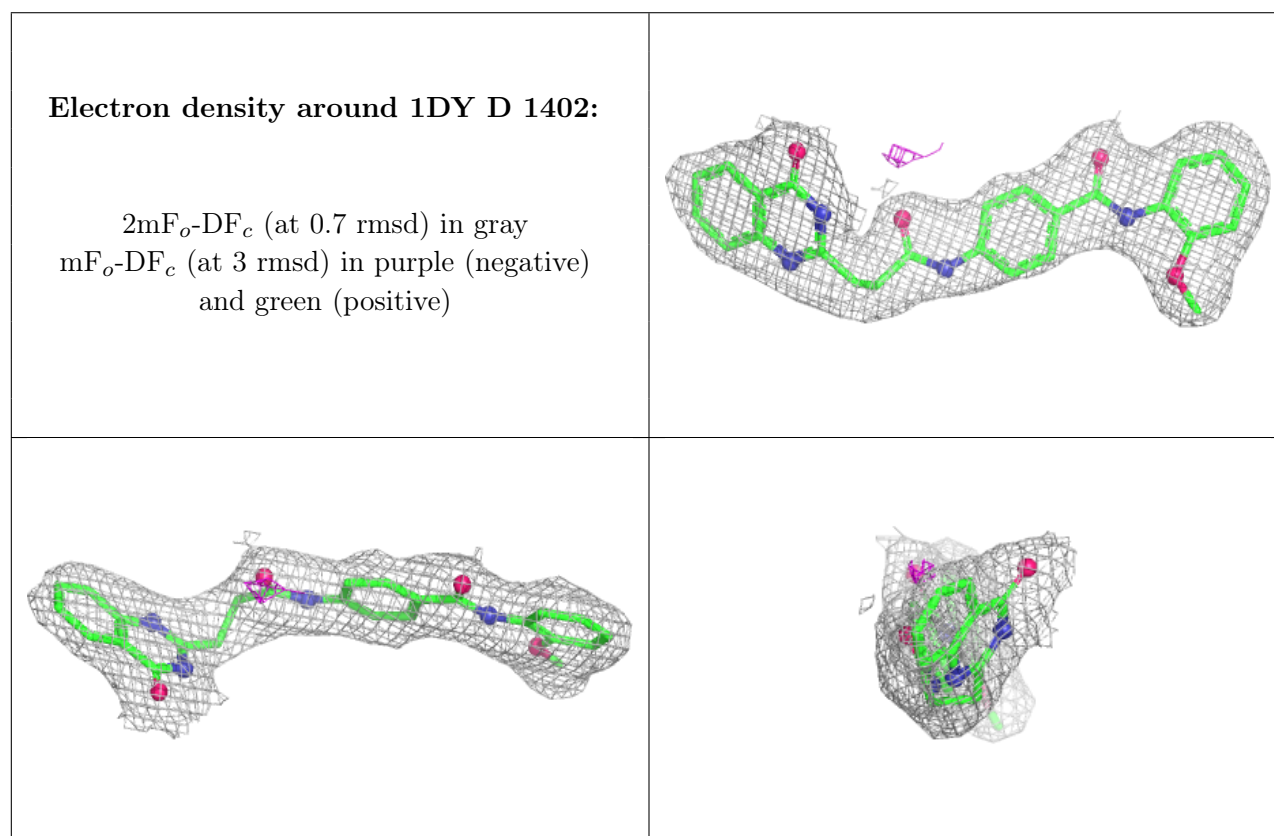
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	1DY	D	1402	33/33	0.93	0.09	25,28,32,34	0
3	1DY	B	1402	33/33	0.95	0.08	21,25,29,30	0
3	1DY	C	1402	33/33	0.95	0.07	23,26,30,31	0
3	1DY	A	1402	33/33	0.95	0.08	24,26,30,33	0
2	ZN	D	1401	1/1	0.97	0.06	66,66,66,66	0
2	ZN	C	1401	1/1	0.98	0.04	59,59,59,59	0
2	ZN	B	1401	1/1	0.98	0.04	48,48,48,48	0

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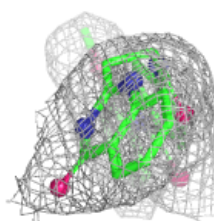
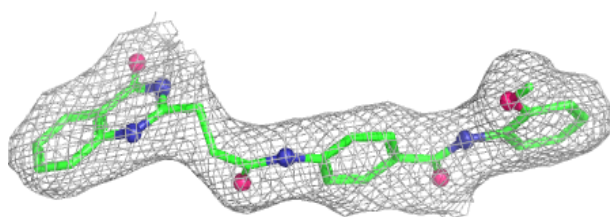
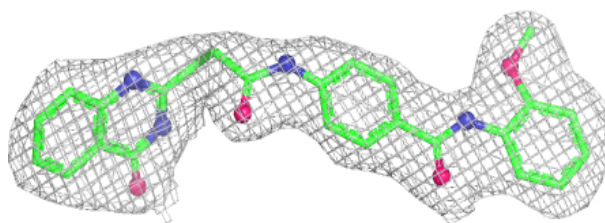
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	A	1401	1/1	0.99	0.02	52,52,52,52	0

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.

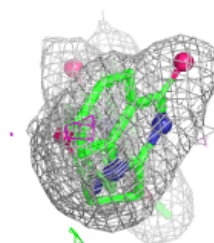
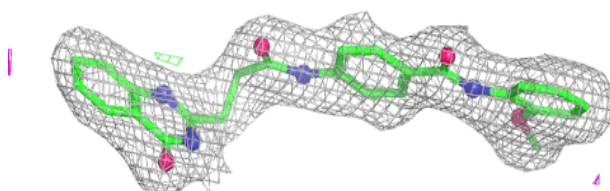
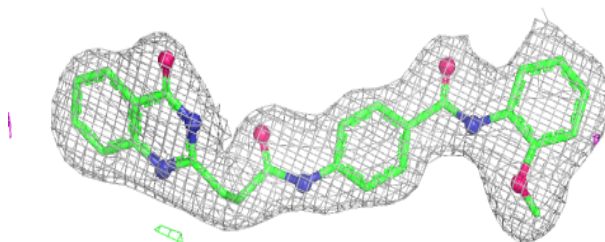


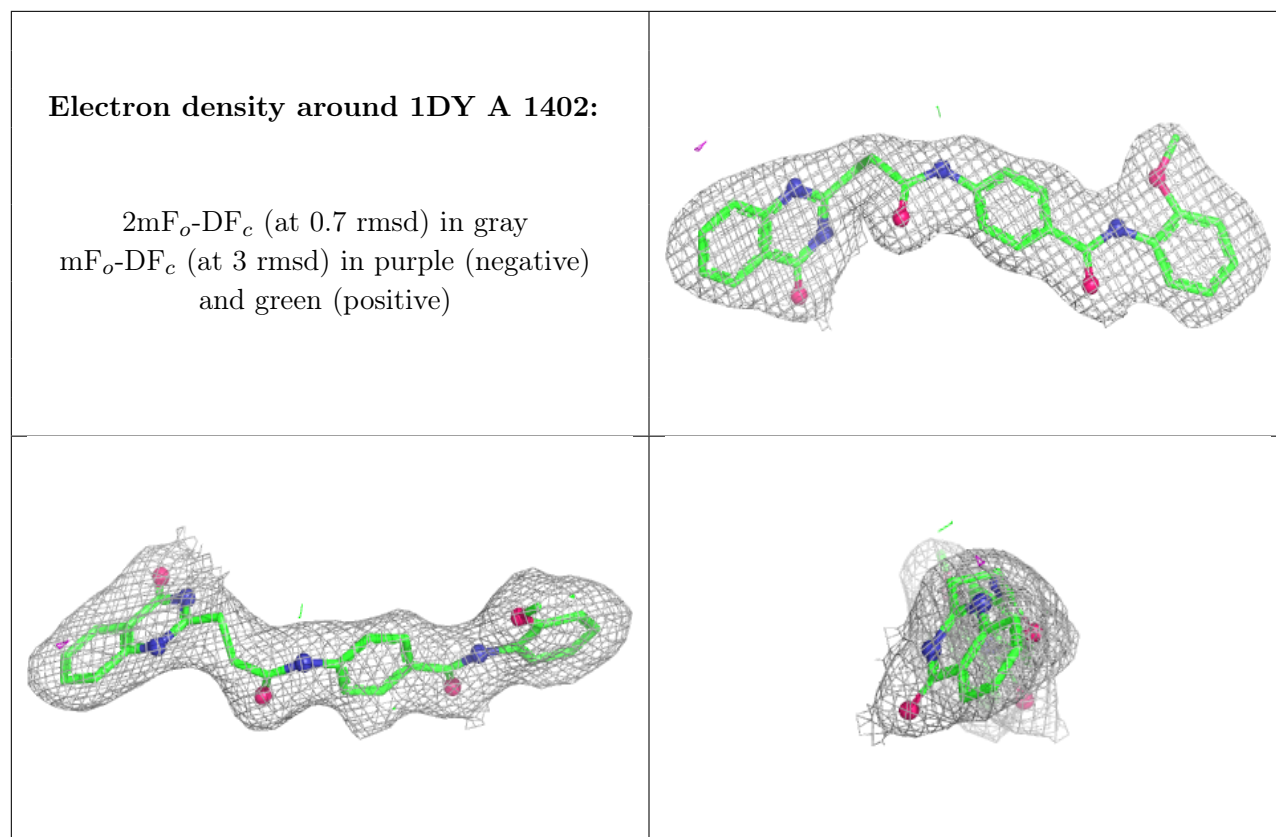
**Electron density around 1DY B 1402:**

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

**Electron density around 1DY C 1402:**

$2mF_o-DF_c$  (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.