



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 16, 2020 – 10:13 AM EDT

PDB ID : 6W2C
Title : Anomalous iodine signal reveals the position of I-paroxetine complexed with the serotonin transporter at the central site
Authors : Coleman, J.A.; Navratna, V.; Yang, D.
Deposited on : 2020-03-05
Resolution : 6.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.4
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) : 2.14.4

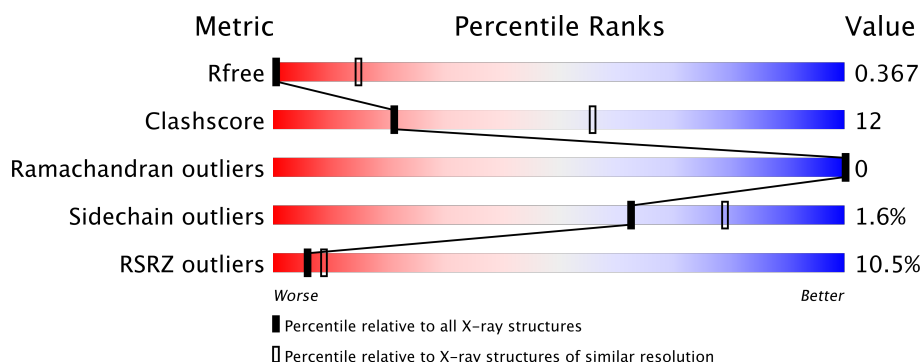
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 6.30 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	1009 (8.70-3.88)
Clashscore	141614	1058 (8.70-3.90)
Ramachandran outliers	138981	1006 (8.70-3.88)
Sidechain outliers	138945	1005 (8.70-3.84)
RSRZ outliers	127900	1018 (8.70-3.78)

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 on 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	549	<div> <div>6%</div> <div>69%</div> <div>28%</div> <div>..</div> </div>
2	B	229	<div> <div>14%</div> <div>71%</div> <div>24%</div> <div>5%</div> </div>
3	C	214	<div> <div>18%</div> <div>75%</div> <div>25%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	RFY	A	705	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7623 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 Sodium-dependent serotonin transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	544	Total	C	N	O	S	0	0	0
			4294	2871	663	736	24			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	GLY	-	expression tag	UNP P31645
A	75	SER	-	expression tag	UNP P31645
A	110	ALA	TYR	conflict	UNP P31645
A	291	ALA	ILE	conflict	UNP P31645
A	554	ALA	CYS	conflict	UNP P31645
A	580	ALA	CYS	conflict	UNP P31645
A	619	LEU	-	expression tag	UNP P31645
A	620	VAL	-	expression tag	UNP P31645
A	621	PRO	-	expression tag	UNP P31645
A	622	ARG	-	expression tag	UNP P31645

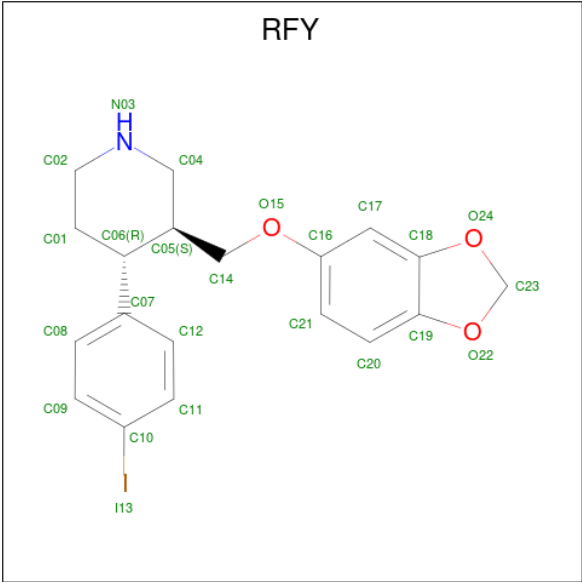
- Molecule 2 is a protein called 8B6 heavy chain antibody fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1643	1038	266	331	8			

- Molecule 3 is a protein called 8B6 light chain antibody fragment.

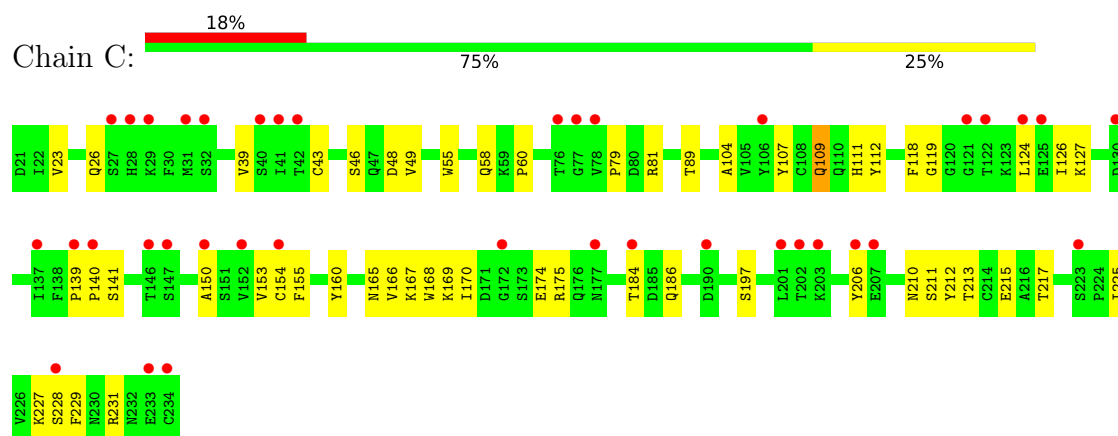
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	0	0	0
			1662	1037	280	337	8			

- Molecule 4 is I-paroxetine (three-letter code: RFY) (formula: C₁₉H₂₀INO₃) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	I	N	O		
4	A	1	24	19	1	1	3	0	0

● Molecule 3: 8B6 light chain antibody fragment



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	126.99Å 159.95Å 140.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 6.30 30.00 – 6.30	Depositor EDS
% Data completeness (in resolution range)	94.6 (30.00-6.30) 92.9 (30.00-6.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.92 (at 6.07Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.306 , 0.369 0.306 , 0.367	Depositor DCC
R_{free} test set	148 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	529.9	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 462.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	7623	wwPDB-VP
Average B, all atoms (Å ²)	568.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: RFY

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	0.37	1/4431 (0.0%)	0.57	4/6054 (0.1%)
2	B	0.33	0/1688	0.52	0/2309
3	C	0.29	0/1700	0.51	0/2307
All	All	0.34	1/7819 (0.0%)	0.55	4/10670 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	81	THR	C-N	6.80	1.49	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	THR	O-C-N	-7.91	110.04	122.70
1	A	81	THR	O-C-N	-7.45	110.78	122.70
1	A	602	GLY	N-CA-C	-5.23	100.03	113.10
1	A	597	LEU	CA-CB-CG	5.05	126.93	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	219	THR	Mainchain

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	4294	0	4263	116	1
2	B	1643	0	1589	33	1
3	C	1662	0	1585	37	0
4	A	24	0	0	7	0
All	All	7623	0	7437	182	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (182) 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:501:VAL:HG21	4:A:705:RFY:I13	1.88	1.43
1:A:594:ALA:O	1:A:598:ILE:HB	1.37	1.20
1:A:497:THR:HB	4:A:705:RFY:I13	2.28	1.02
1:A:501:VAL:CG2	4:A:705:RFY:I13	2.82	0.97
3:C:211:SER:HA	3:C:229:PHE:O	1.69	0.93
1:A:172:ILE:HD11	4:A:705:RFY:I13	2.49	0.82
2:B:30:LEU:HD22	2:B:172:PRO:HD3	1.63	0.80
1:A:526:LYS:HD3	1:A:532:SER:HB2	1.64	0.80
3:C:126:ILE:O	3:C:186:GLN:NE2	2.14	0.79
3:C:26:GLN:HE21	3:C:119:GLY:HA3	1.51	0.76
1:A:461:ARG:HH11	1:A:461:ARG:HG3	1.53	0.73
1:A:137:LEU:HD12	1:A:348:ALA:HB2	1.69	0.73
2:B:192:PRO:HD3	3:C:184:THR:HG22	1.72	0.71
1:A:141:GLN:HE21	1:A:351:ASN:HB3	1.55	0.70
1:A:307:ARG:NE	1:A:388:GLU:OE2	2.23	0.70
1:A:139:LEU:HD11	1:A:512:VAL:HG11	1.72	0.69
3:C:213:THR:HA	3:C:227:LYS:O	1.93	0.67
3:C:210:ASN:OD1	3:C:231:ARG:N	2.27	0.67
3:C:23:VAL:H	3:C:46:SER:HB3	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:215:GLU:HA	3:C:225:ILE:O	1.96	0.66
1:A:307:ARG:HD3	1:A:392:GLU:HG3	1.77	0.66
1:A:136:GLU:HG2	1:A:344:LEU:HB2	1.77	0.65
2:B:174:SER:HB3	2:B:225:PRO:HG2	1.79	0.65
1:A:437:ASP:N	1:A:437:ASP:OD1	2.30	0.64
2:B:187:SER:OG	2:B:207:THR:O	2.11	0.64
1:A:206:THR:HG23	1:A:234:ARG:NH2	2.13	0.64
1:A:87:ASP:OD1	1:A:282:TRP:NE1	2.26	0.64
1:A:363:VAL:O	1:A:367:VAL:HG23	1.98	0.63
1:A:447:ILE:HA	1:A:465:PHE:HE2	1.63	0.63
2:B:163:LEU:HB3	2:B:235:LEU:HD22	1.81	0.62
1:A:98:ASP:HB2	1:A:176:TYR:OH	2.01	0.61
3:C:165:ASN:HB3	3:C:217:THR:HB	1.82	0.61
1:A:101:ASN:ND2	1:A:372:SER:OG	2.33	0.61
1:A:184:LEU:HD21	1:A:261:LEU:HD23	1.82	0.60
1:A:157:ILE:HG23	1:A:611:SER:HB3	1.84	0.60
1:A:184:LEU:HA	1:A:432:ILE:HD11	1.85	0.59
3:C:168:TRP:O	3:C:174:GLU:HA	2.04	0.58
1:A:292:LEU:O	1:A:296:LEU:HB2	2.04	0.58
1:A:327:ILE:HD13	1:A:555:SER:CB	2.34	0.58
3:C:206:TYR:CZ	3:C:231:ARG:HG3	2.39	0.58
1:A:352:LYS:HB3	1:A:355:ASN:HB2	1.85	0.57
2:B:86:LYS:NZ	2:B:104:SER:O	2.38	0.57
2:B:149:LEU:HD13	3:C:153:VAL:HG21	1.86	0.57
1:A:461:ARG:NH1	1:A:461:ARG:HG3	2.20	0.56
3:C:212:TYR:HB2	3:C:229:PHE:CE1	2.40	0.56
3:C:139:PRO:HB3	3:C:229:PHE:CE2	2.41	0.56
3:C:212:TYR:O	3:C:228:SER:HA	2.06	0.56
1:A:599:ILE:O	1:A:601:PRO:HD3	2.06	0.56
1:A:95:TYR:O	4:A:705:RFY:N03	2.39	0.56
1:A:453:GLU:HG3	1:A:454:PHE:CD1	2.42	0.55
1:A:478:LEU:HD23	1:A:481:LEU:HD12	1.88	0.55
2:B:139:ALA:HB3	2:B:171:PHE:CE2	2.42	0.55
2:B:148:PRO:O	3:C:141:SER:HB3	2.07	0.55
1:A:125:ALA:HA	1:A:333:ILE:HG12	1.89	0.55
2:B:149:LEU:HD11	2:B:166:LEU:HB2	1.89	0.55
3:C:169:LYS:HB2	3:C:213:THR:O	2.06	0.54
1:A:312:TYR:CZ	1:A:376:GLY:HA3	2.42	0.54
1:A:442:GLY:HA3	4:A:705:RFY:C18	2.38	0.54
1:A:87:ASP:O	1:A:91:SER:HB2	2.08	0.54
2:B:52:TYR:CE2	2:B:120:VAL:HG12	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:ILE:HA	1:A:440:PHE:HE1	1.73	0.53
1:A:95:TYR:CD1	1:A:343:VAL:HG11	2.43	0.53
1:A:393:ASP:O	1:A:395:SER:N	2.42	0.53
3:C:48:ASP:HA	3:C:89:THR:HG22	1.91	0.53
1:A:485:GLY:O	1:A:489:VAL:HG23	2.09	0.52
1:A:603:THR:HG22	1:A:604:PHE:H	1.75	0.52
1:A:136:GLU:CG	1:A:344:LEU:HB2	2.40	0.52
3:C:231:ARG:O	3:C:231:ARG:HG2	2.10	0.52
1:A:79:ARG:NE	1:A:349:SER:OG	2.44	0.51
1:A:392:GLU:HG2	1:A:393:ASP:H	1.75	0.51
2:B:180:ASN:HD21	2:B:218:VAL:HG13	1.75	0.51
3:C:111:HIS:O	3:C:111:HIS:ND1	2.43	0.51
1:A:125:ALA:O	1:A:130:ILE:HG12	2.10	0.51
1:A:322:GLU:O	1:A:325:VAL:HG22	2.11	0.51
2:B:54:ASN:ND2	2:B:118:SER:OG	2.35	0.50
1:A:234:ARG:HE	1:A:240:HIS:CG	2.29	0.50
2:B:163:LEU:O	2:B:205:SER:HA	2.12	0.50
2:B:59:SER:HB2	2:B:62:LYS:HB2	1.93	0.50
1:A:327:ILE:HD13	1:A:555:SER:HB2	1.93	0.50
1:A:175:TYR:HB2	1:A:489:VAL:HG13	1.93	0.50
1:A:195:LEU:H	1:A:195:LEU:HD12	1.77	0.49
1:A:106:PRO:HG3	1:A:376:GLY:HA2	1.94	0.49
1:A:113:GLY:HA2	1:A:316:ASN:HB3	1.93	0.49
3:C:43:CYS:HB2	3:C:55:TRP:CH2	2.47	0.49
1:A:518:ILE:HD11	1:A:541:TRP:CE3	2.47	0.49
1:A:311:PHE:CE1	1:A:395:SER:HB2	2.47	0.49
1:A:454:PHE:O	1:A:458:TRP:HB2	2.13	0.49
1:A:272:LYS:HD3	1:A:276:THR:HG23	1.95	0.48
1:A:400:ASP:OD1	1:A:400:ASP:N	2.47	0.48
1:A:312:TYR:OH	1:A:376:GLY:HA3	2.14	0.48
1:A:263:PHE:HA	1:A:266:ILE:HG12	1.96	0.48
1:A:603:THR:H	1:A:606:GLU:HB3	1.77	0.48
2:B:170:TYR:OH	2:B:175:VAL:HG11	2.14	0.48
3:C:167:LYS:HD2	3:C:215:GLU:OE1	2.14	0.48
3:C:140:PRO:HG2	3:C:150:ALA:HB1	1.95	0.47
3:C:170:ILE:HG13	3:C:175:ARG:HB2	1.95	0.47
2:B:205:SER:HB3	3:C:155:PHE:CE2	2.49	0.47
2:B:48:PHE:CE2	2:B:72:PRO:HB3	2.50	0.47
2:B:183:SER:HA	2:B:184:LEU:HA	1.58	0.47
2:B:58:GLN:O	2:B:111:ALA:HB1	2.14	0.47
2:B:135:THR:HG21	2:B:172:PRO:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:VAL:O	1:A:255:LEU:HD21	2.15	0.46
1:A:256:ALA:O	1:A:260:MET:HB2	2.13	0.46
1:A:271:TRP:HA	1:A:466:VAL:HG11	1.97	0.46
1:A:206:THR:HG23	1:A:234:ARG:HH22	1.78	0.46
1:A:263:PHE:O	1:A:267:TYR:HB2	2.15	0.46
1:A:156:PRO:HG2	1:A:611:SER:O	2.15	0.46
1:A:296:LEU:HB2	1:A:378:VAL:HG22	1.97	0.46
3:C:166:VAL:HG11	3:C:197:SER:CB	2.45	0.46
3:C:39:VAL:HG21	3:C:124:LEU:HD11	1.98	0.46
1:A:136:GLU:HG2	1:A:344:LEU:HD12	1.98	0.46
1:A:178:THR:HG21	1:A:480:THR:HB	1.98	0.45
1:A:82:TRP:HD1	1:A:83:GLY:H	1.62	0.45
1:A:103:TRP:C	1:A:106:PRO:HD2	2.37	0.45
1:A:91:SER:HA	1:A:281:VAL:HG11	1.98	0.45
2:B:164:GLY:HA2	2:B:204:SER:O	2.16	0.45
1:A:455:PRO:O	1:A:457:VAL:N	2.47	0.45
1:A:407:PHE:HD1	1:A:428:PHE:HE1	1.65	0.45
1:A:545:SER:HB2	1:A:546:PRO:HD3	1.99	0.45
1:A:148:ILE:HG13	1:A:449:ALA:HB1	1.98	0.45
3:C:127:LYS:HA	3:C:160:TYR:OH	2.17	0.45
3:C:206:TYR:O	3:C:231:ARG:HD3	2.17	0.45
3:C:154:CYS:HB2	3:C:168:TRP:CH2	2.52	0.45
1:A:142:TYR:HD2	1:A:143:HIS:CE1	2.35	0.44
1:A:311:PHE:CZ	1:A:395:SER:HB2	2.52	0.44
2:B:58:GLN:NE2	2:B:114:TYR:HE2	2.14	0.44
2:B:160:SER:HA	2:B:209:PRO:HA	2.00	0.44
1:A:178:THR:HG22	1:A:259:ILE:HD12	1.99	0.44
1:A:390:ARG:CZ	1:A:397:VAL:HG23	2.47	0.44
2:B:53:MET:HG2	2:B:98:ALA:CB	2.46	0.44
1:A:439:THR:HA	4:A:705:RFY:C20	2.47	0.44
1:A:343:VAL:CG1	1:A:441:ALA:HB1	2.48	0.44
1:A:396:GLU:HA	1:A:399:LYS:HG3	2.00	0.44
3:C:168:TRP:NE1	3:C:197:SER:OG	2.36	0.44
1:A:270:ILE:HD13	1:A:277:SER:HB3	2.00	0.43
1:A:589:ILE:HG23	1:A:590:PRO:HD3	1.99	0.43
1:A:105:PHE:HA	1:A:108:ILE:HG22	2.00	0.43
1:A:102:VAL:HG21	1:A:434:LEU:CD1	2.48	0.43
2:B:218:VAL:O	2:B:234:LYS:HG3	2.19	0.43
2:B:72:PRO:HA	2:B:91:VAL:HG11	2.01	0.43
1:A:284:THR:OG1	1:A:437:ASP:HB3	2.18	0.43
1:A:429:LEU:O	1:A:433:THR:OG1	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:VAL:HG21	1:A:434:LEU:HD13	2.00	0.42
1:A:599:ILE:HG23	1:A:600:THR:HG23	2.01	0.42
1:A:221:THR:HG22	1:A:223:HIS:H	1.85	0.42
2:B:55:TRP:CD1	2:B:89:LEU:HD22	2.55	0.42
1:A:103:TRP:CH2	1:A:402:GLY:HA2	2.54	0.42
1:A:263:PHE:HD1	1:A:266:ILE:HD11	1.84	0.42
1:A:302:LEU:HD13	1:A:385:TYR:CD2	2.54	0.42
1:A:352:LYS:HE2	1:A:355:ASN:HB2	2.02	0.42
1:A:453:GLU:HG3	1:A:454:PHE:CE1	2.55	0.42
1:A:563:LEU:HD12	1:A:563:LEU:H	1.84	0.42
1:A:302:LEU:HD13	1:A:385:TYR:CG	2.55	0.42
1:A:577:LEU:O	1:A:581:ILE:HG13	2.20	0.42
1:A:180:MET:HE1	1:A:436:LEU:HG	2.02	0.41
1:A:103:TRP:HD1	1:A:406:LEU:HD11	1.85	0.41
2:B:189:HIS:HB2	2:B:205:SER:OG	2.20	0.41
2:B:69:ASN:OD1	2:B:78:SER:HB3	2.20	0.41
1:A:159:LYS:HB3	1:A:592:TYR:CE2	2.55	0.41
3:C:109:GLN:HB2	3:C:118:PHE:CD2	2.55	0.41
1:A:275:LYS:O	1:A:279:LYS:HD3	2.20	0.41
1:A:608:ILE:O	1:A:612:ILE:HG23	2.19	0.41
3:C:127:LYS:HD2	3:C:160:TYR:CE2	2.55	0.41
3:C:58:GLN:HE21	3:C:107:TYR:HE1	1.68	0.41
3:C:49:VAL:HG13	3:C:112:TYR:CD1	2.56	0.41
3:C:60:PRO:HD3	3:C:104:ALA:HA	2.03	0.41
1:A:122:THR:O	1:A:125:ALA:HB3	2.21	0.41
1:A:392:GLU:CG	1:A:393:ASP:H	2.34	0.41
3:C:79:PRO:HB2	3:C:81:ARG:HG2	2.03	0.41
1:A:563:LEU:HD13	1:A:570:TYR:CG	2.56	0.41
1:A:518:ILE:HA	1:A:518:ILE:HD12	1.82	0.40
2:B:31:VAL:HG21	2:B:105:LEU:HD13	2.03	0.40
1:A:298:ARG:NH2	1:A:301:THR:OG1	2.54	0.40
2:B:188:VAL:HG22	2:B:206:VAL:HG23	2.03	0.40
1:A:103:TRP:CD1	1:A:406:LEU:HD11	2.56	0.40
1:A:287:PHE:HB3	1:A:433:THR:HG21	2.04	0.40
1:A:453:GLU:HG3	1:A:454:PHE:HD1	1.84	0.40
2:B:57:LYS:HB2	2:B:67:ILE:HD11	2.03	0.40
1:A:142:TYR:HD2	1:A:143:HIS:ND1	2.20	0.40
1:A:195:LEU:HA	1:A:196:PRO:HD3	1.98	0.40
1:A:426:ILE:O	1:A:430:MET:HB2	2.21	0.40
1:A:232:TYR:HA	1:A:236:VAL:HG12	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:609:ILE:CD1	2:B:216:GLN:NE2[4_577]	2.11	0.09

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	542/549 (99%)	519 (96%)	23 (4%)	0	100	100
2	B	216/229 (94%)	207 (96%)	9 (4%)	0	100	100
3	C	212/214 (99%)	200 (94%)	12 (6%)	0	100	100
All	All	970/992 (98%)	926 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	450/462 (97%)	438 (97%)	12 (3%)	44	65
2	B	190/201 (94%)	190 (100%)	0	100	100
3	C	189/190 (100%)	188 (100%)	1 (0%)	88	93
All	All	829/853 (97%)	816 (98%)	13 (2%)	62	79

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ARG
1	A	82	TRP
1	A	217	ASN
1	A	260	MET
1	A	271	TRP
1	A	437	ASP
1	A	460	LYS
1	A	461	ARG
1	A	494	GLU
1	A	563	LEU
1	A	604	PHE
1	A	605	LYS
3	C	109	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	141	GLN
1	A	217	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](#)

1 ligand is 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
4	RFY	A	705	-	26,27,27	1.54	5 (19%)	33,37,37	1.48	7 (21%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	RFY	A	705	-	-	5/9/26/26	1/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	705	RFY	C07-C06	4.11	1.58	1.51
4	A	705	RFY	C01-C06	-2.71	1.50	1.53
4	A	705	RFY	O24-C18	2.58	1.42	1.38
4	A	705	RFY	O22-C19	2.55	1.42	1.38
4	A	705	RFY	C14-C05	2.04	1.55	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	705	RFY	C01-C06-C05	3.39	113.55	109.62
4	A	705	RFY	C07-C06-C05	-3.05	107.76	113.69
4	A	705	RFY	O24-C18-C17	3.02	131.89	127.85
4	A	705	RFY	C14-C05-C06	-2.33	107.85	112.05
4	A	705	RFY	C01-C06-C07	-2.23	108.48	112.57
4	A	705	RFY	O22-C19-C20	2.23	131.86	127.81
4	A	705	RFY	C14-O15-C16	-2.13	113.41	117.93

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	705	RFY	C17-C16-O15-C14
4	A	705	RFY	C21-C16-O15-C14

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Mol	Chain	Res	Type	Atoms
4	A	705	RFY	C05-C14-O15-C16
4	A	705	RFY	C06-C05-C14-O15
4	A	705	RFY	C04-C05-C14-O15

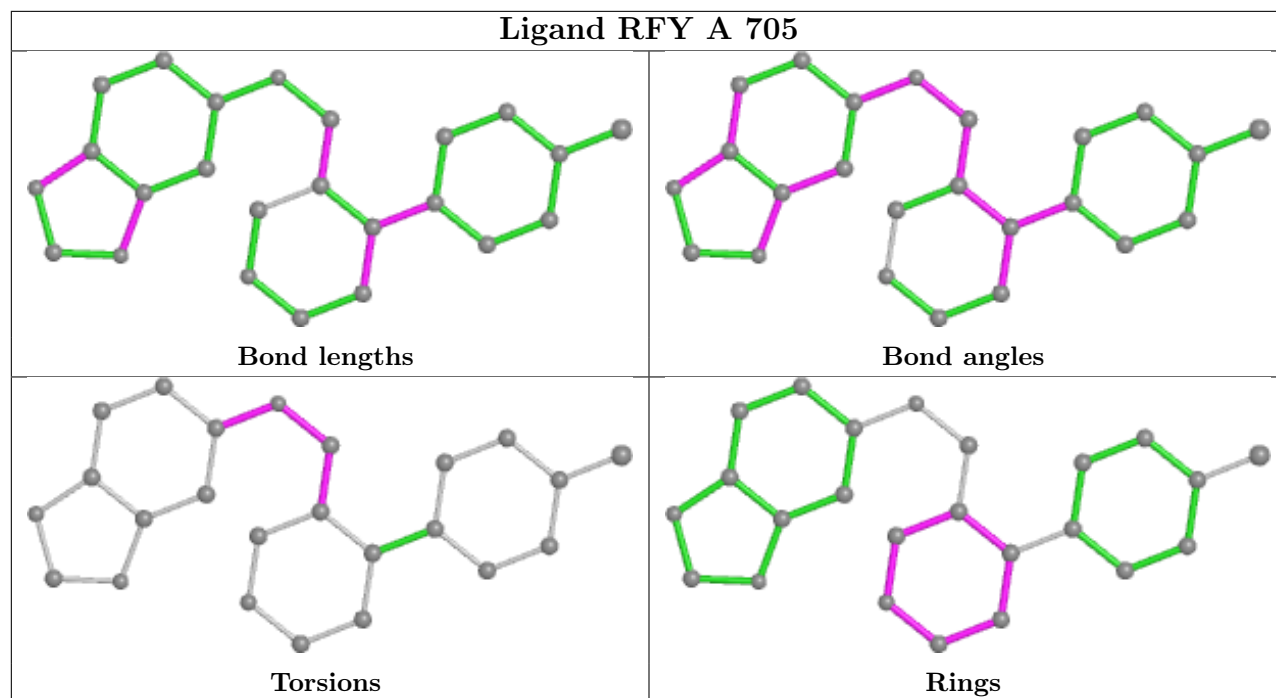
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	705	RFY	C01-C02-C04-C05-C06-N03

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	705	RFY	7	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, 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	544/549 (99%)	0.26	32 (5%)	22 22	541, 553, 577, 612	0
2	B	218/229 (95%)	0.89	32 (14%)	2 5	536, 559, 643, 665	0
3	C	214/214 (100%)	0.95	38 (17%)	1 4	547, 581, 641, 662	0
All	All	976/992 (98%)	0.55	102 (10%)	6 9	536, 556, 635, 665	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	SER	8.9
1	A	74	GLY	7.8
1	A	532	SER	7.6
1	A	76	GLN	6.0
1	A	531	PHE	5.9
1	A	338	GLY	5.8
3	C	150	ALA	5.6
2	B	153	CYS	5.5
1	A	339	PRO	5.4
3	C	140	PRO	5.3
3	C	122	THR	5.3
2	B	154	GLY	5.3
1	A	533	PRO	5.1
1	A	340	GLY	4.9
3	C	234	CYS	4.9
2	B	157	THR	4.8
1	A	77	GLY	4.6
2	B	209	PRO	4.5
3	C	154	CYS	4.3
2	B	211	SER	4.2
2	B	144	PRO	4.2
2	B	212	THR	4.2
3	C	139	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
2	B	158	GLY	4.1
3	C	121	GLY	4.0
2	B	156	THR	3.9
1	A	411	ALA	3.7
2	B	210	SER	3.7
3	C	78	VAL	3.6
3	C	28	HIS	3.6
2	B	237	PRO	3.6
3	C	41	ILE	3.6
3	C	233	GLU	3.5
1	A	530	GLY	3.5
3	C	203	LYS	3.4
1	A	134	TYR	3.3
3	C	202	THR	3.2
1	A	529	LEU	3.1
2	B	213	TRP	3.1
2	B	208	VAL	3.1
3	C	77	GLY	3.1
1	A	101	ASN	3.0
2	B	91	VAL	2.9
1	A	317	TRP	2.9
2	B	226	ALA	2.9
3	C	146	THR	2.9
3	C	152	VAL	2.9
3	C	172	GLY	2.8
2	B	103	ARG	2.7
3	C	29	LYS	2.7
1	A	407	PHE	2.7
3	C	125	GLU	2.6
1	A	342	GLY	2.6
2	B	85	GLY	2.6
2	B	28	PRO	2.6
3	C	206	TYR	2.5
3	C	27	SER	2.5
2	B	36	SER	2.5
3	C	130	ASP	2.5
1	A	183	ALA	2.5
1	A	605	LYS	2.4
3	C	42	THR	2.4
1	A	341	PHE	2.4
3	C	201	LEU	2.4
3	C	124	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	159	SER	2.4
2	B	167	VAL	2.3
3	C	76	THR	2.3
2	B	37	VAL	2.3
3	C	32	SER	2.3
1	A	561	PRO	2.3
1	A	534	GLY	2.3
1	A	358	TYR	2.2
1	A	142	TYR	2.2
2	B	172	PRO	2.2
3	C	177	ASN	2.2
2	B	203	SER	2.2
1	A	320	LEU	2.2
3	C	137	ILE	2.2
2	B	236	GLU	2.2
1	A	428	PHE	2.2
3	C	190	ASP	2.2
3	C	184	THR	2.2
3	C	223	SER	2.2
2	B	192	PRO	2.2
2	B	228	SER	2.2
3	C	228	SER	2.2
1	A	136	GLU	2.1
3	C	31	MET	2.1
2	B	43	ALA	2.1
3	C	40	SER	2.1
2	B	102	LEU	2.1
1	A	79	ARG	2.1
2	B	92	ASP	2.1
2	B	141	THR	2.1
3	C	106	TYR	2.1
1	A	87	ASP	2.0
1	A	178	THR	2.0
2	B	164	GLY	2.0
3	C	147	SER	2.0
1	A	465	PHE	2.0
3	C	207	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

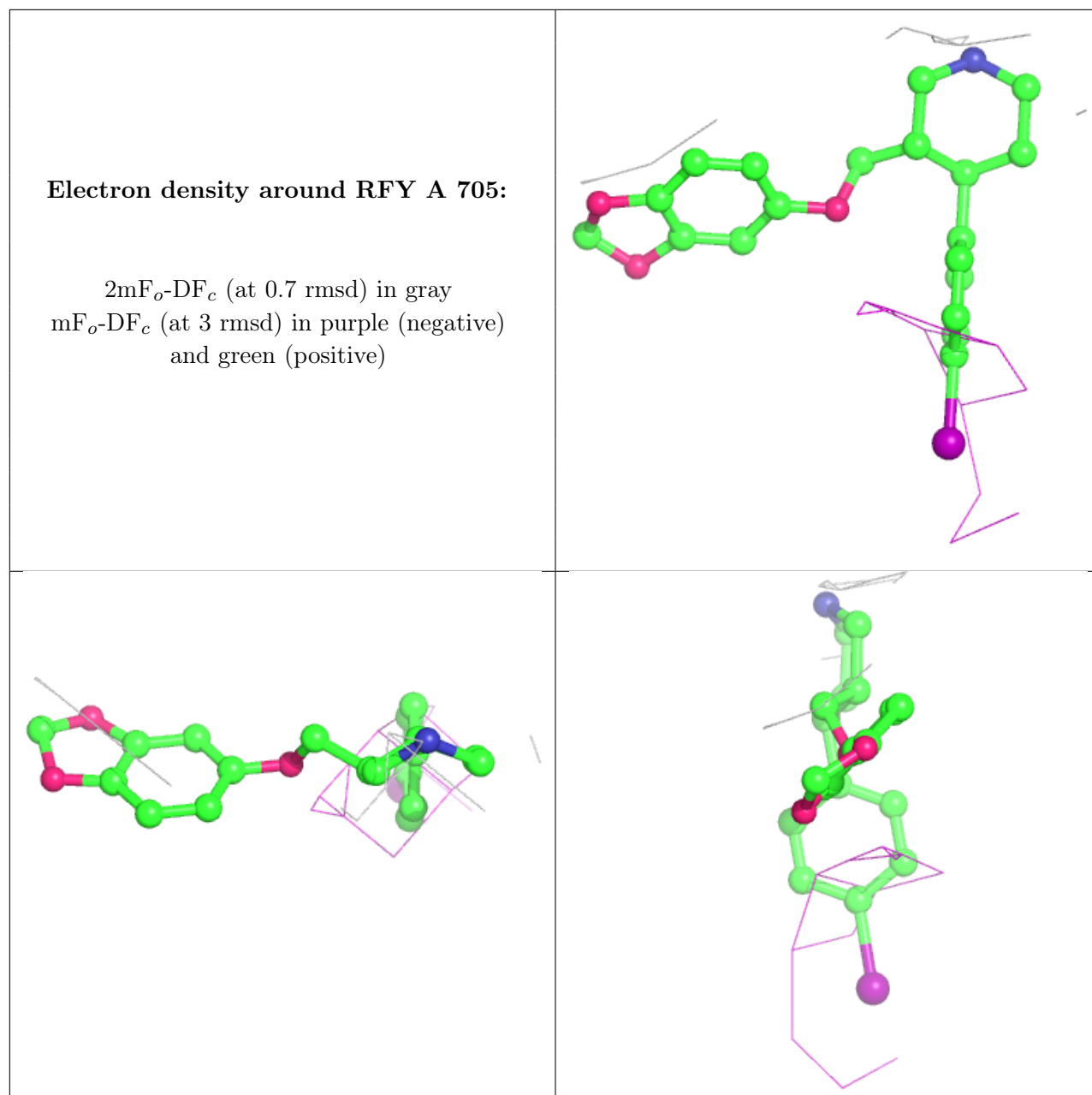
There are no monosaccharides in this entry.

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, 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(\AA^2)	Q<0.9
4	RFY	A	705	24/24	0.45	0.62	539,542,545,552	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.