Full wwPDB X-ray Structure Validation Report ⓘ

Dec 10, 2023 – 03:33 pm GMT

PDB ID	:	2V91
Title	:	STRUCTURE OF STRICTOSIDINE SYNTHASE IN COMPLEX WITH STRICTOSIDINE
Authors	:	Loris, E.A.; Panjikar, S.; Ruppert, M.; Barleben, L.; Unger, M.; Stoeckigt, J.
Deposited on	:	2007-08-16
Resolution	:	3.01 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

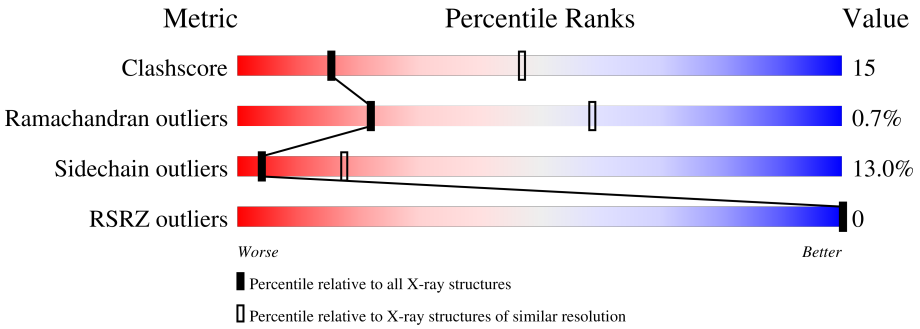
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.01 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))
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)
RSRZ outliers	127900	2287 (3.04-3.00)

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 <=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	302	67% 25% 7%
1	B	302	66% 30% .

2 Entry composition

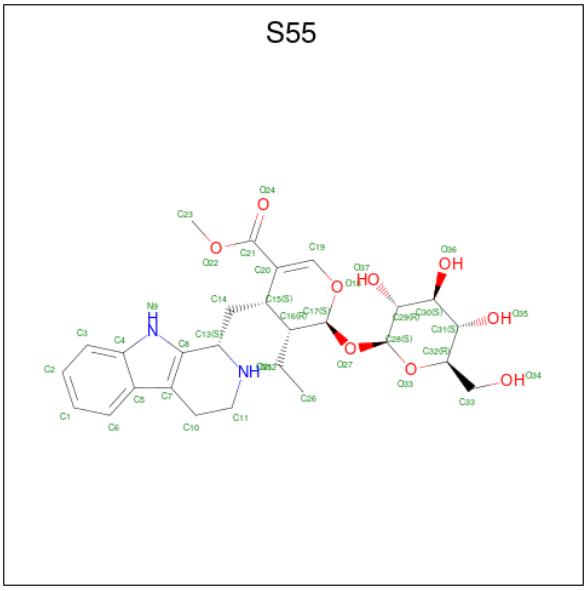
There are 3 unique types of molecules in this entry. The entry contains 4858 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 STRICTOSIDINE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2363	1511	388	459	5			
1	B	302	Total	C	N	O	S	0	0	0
			2364	1512	388	459	5			

- Molecule 2 is METHYL (2S,3R,4S)-3-ETHYL-2-(BETA-D-GLUCOPYRANOSYLOXY)-4-[(1S)-2,3,4,9-TETRAHYDRO-1H-BETA-CARBOLIN-1-YLMETHYL]-3,4-DIHYDRO-2H-PYRAN-5-CARBOXYLATE (three-letter code: S55) (formula: $C_{27}H_{36}N_2O_9$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			38	27	2	9		
2	B	1	Total	C	N	O	0	0
			38	27	2	9		

- Molecule 3 is water.

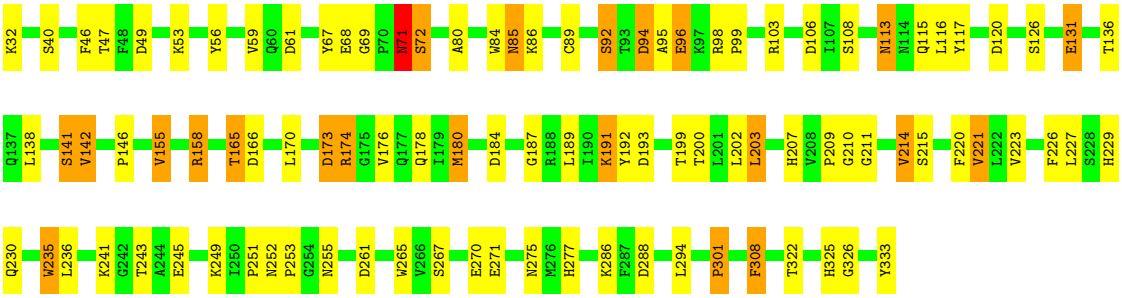
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	35	Total O 35 35	0	0
3	B	20	Total O 20 20	0	0

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.

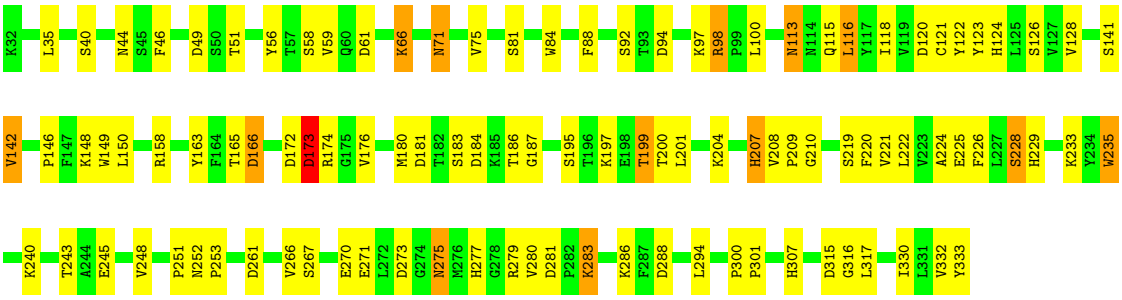
• Molecule 1: STRICTOSIDINE SYNTHASE

Chain A:



• Molecule 1: STRICTOSIDINE SYNTHASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	150.18Å 150.18Å 121.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.01 19.77 – 3.01	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-3.01) 100.0 (19.77-3.01)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.47 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.192 , 0.235 0.182 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	60.4	Xtriage
Anisotropy	0.758	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 31.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.023 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4858	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: S55

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	0.89	5/2428 (0.2%)	1.00	10/3309 (0.3%)
1	B	0.79	0/2429	0.99	11/3311 (0.3%)
All	All	0.84	5/4857 (0.1%)	0.99	21/6620 (0.3%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	333	TYR	CB-CG	6.30	1.61	1.51
1	A	333	TYR	CD1-CE1	6.30	1.48	1.39
1	A	333	TYR	CD2-CE2	6.08	1.48	1.39
1	A	180	MET	SD-CE	5.46	2.08	1.77
1	A	71	ASN	CB-CG	5.30	1.63	1.51

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	166	ASP	CB-CG-OD2	7.52	125.07	118.30
1	B	288	ASP	CB-CG-OD2	6.81	124.43	118.30
1	B	172	ASP	CB-CG-OD2	6.50	124.14	118.30
1	A	106	ASP	CB-CG-OD2	6.42	124.08	118.30
1	B	120	ASP	CB-CG-OD2	6.37	124.04	118.30
1	A	120	ASP	CB-CG-OD2	6.35	124.01	118.30
1	B	261	ASP	CB-CG-OD2	6.29	123.96	118.30
1	A	94	ASP	CB-CG-OD2	6.13	123.82	118.30
1	B	184	ASP	CB-CG-OD2	5.97	123.67	118.30
1	A	61	ASP	CB-CG-OD1	5.68	123.41	118.30
1	B	273	ASP	CB-CG-OD2	5.66	123.39	118.30
1	A	184	ASP	CB-CG-OD2	5.64	123.38	118.30
1	B	315	ASP	CB-CG-OD2	5.49	123.24	118.30
1	B	94	ASP	CB-CG-OD2	5.33	123.10	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	49	ASP	CB-CG-OD1	5.25	123.03	118.30
1	B	166	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	261	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	180	MET	CG-SD-CE	-5.13	91.99	100.20
1	B	98	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	193	ASP	CB-CG-OD2	5.11	122.90	118.30
1	A	120	ASP	CB-CG-OD1	-5.06	113.74	118.30

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	2363	0	2201	68	0
1	B	2364	0	2202	62	0
2	A	38	0	36	7	0
2	B	38	0	36	4	0
3	A	35	0	0	3	0
3	B	20	0	0	0	0
All	All	4858	0	4475	137	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 15.

All (137) 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:180:MET:CE	1:A:180:MET:SD	2.08	1.41
2:A:1334:S55:H263	2:A:1334:S55:H13	1.26	1.13
1:B:229:HIS:CD2	1:B:252:ASN:H	1.91	0.88
2:B:1334:S55:H263	2:B:1334:S55:H13	1.63	0.80
1:B:141:SER:HB3	1:B:146:PRO:HA	1.65	0.79
1:B:229:HIS:HD2	1:B:251:PRO:HA	1.49	0.77
1:A:229:HIS:CD2	1:A:252:ASN:H	2.03	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:VAL:HG22	1:A:236:LEU:HD11	1.68	0.76
1:B:142:VAL:HG23	1:B:199:THR:HB	1.69	0.72
1:B:252:ASN:ND2	1:B:271:GLU:H	1.87	0.72
1:B:124:HIS:HA	1:B:150:LEU:HD12	1.73	0.71
1:A:116:LEU:HD12	1:A:117:TYR:N	2.05	0.70
1:A:131:GLU:H	1:A:131:GLU:CD	1.97	0.69
1:B:229:HIS:HD2	1:B:252:ASN:H	1.38	0.69
1:A:141:SER:HB2	1:A:146:PRO:HA	1.74	0.68
1:B:229:HIS:CD2	1:B:251:PRO:HA	2.29	0.67
1:A:308:PHE:O	2:A:1334:S55:H112	1.93	0.67
1:A:252:ASN:ND2	1:A:271:GLU:H	1.93	0.65
1:B:280:VAL:HG21	1:B:307:HIS:CE1	2.32	0.64
1:A:80:ALA:HB1	3:A:2018:HOH:O	1.96	0.64
1:A:191:LYS:HD3	1:A:200:THR:HB	1.80	0.64
1:A:288:ASP:HB3	1:A:294:LEU:HD11	1.80	0.64
1:A:158:ARG:HH11	1:A:158:ARG:CB	2.11	0.63
1:A:71:ASN:HD22	1:A:72:SER:HB2	1.63	0.62
1:A:158:ARG:HH11	1:A:158:ARG:HB2	1.64	0.62
1:B:252:ASN:HD22	1:B:270:GLU:HA	1.66	0.61
2:A:1334:S55:H15	2:A:1334:S55:H9	1.66	0.60
1:A:165:THR:HG21	1:A:210:GLY:O	2.01	0.60
1:A:230:GLN:HE21	1:A:249:LYS:HE3	1.67	0.60
1:A:180:MET:HB3	1:A:180:MET:HE3	1.81	0.60
1:A:142:VAL:HG23	1:A:199:THR:HG22	1.83	0.60
2:A:1334:S55:H263	2:A:1334:S55:C13	2.18	0.59
1:B:229:HIS:HD2	1:B:252:ASN:N	2.01	0.59
1:A:158:ARG:HB2	1:A:158:ARG:NH1	2.18	0.57
1:A:326:GLY:H	1:B:71:ASN:HA	1.69	0.57
1:A:243:THR:HG21	3:A:2031:HOH:O	2.05	0.57
1:A:275:ASN:ND2	1:A:277:HIS:H	2.02	0.56
1:B:316:GLY:O	1:B:317:LEU:HD23	2.05	0.56
1:B:240:LYS:O	1:B:243:THR:HB	2.05	0.56
1:A:113:ASN:ND2	1:A:115:GLN:HG3	2.20	0.56
1:B:208:VAL:O	1:B:208:VAL:HG12	2.05	0.56
1:A:180:MET:CE	1:A:180:MET:HB3	2.36	0.55
1:A:180:MET:CE	1:A:180:MET:CG	2.84	0.55
1:A:71:ASN:ND2	1:A:72:SER:HB2	2.22	0.54
1:A:187:GLY:HA3	1:A:209:PRO:CD	2.38	0.54
1:A:252:ASN:HD22	1:A:270:GLU:HA	1.72	0.54
1:A:202:LEU:O	1:A:203:LEU:HD23	2.08	0.53
2:A:1334:S55:H15	2:A:1334:S55:N9	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:235:TRP:HZ3	1:B:245:GLU:CG	2.22	0.53
1:A:275:ASN:HD22	1:A:277:HIS:H	1.55	0.53
2:A:1334:S55:H13	2:A:1334:S55:C26	2.19	0.53
1:B:141:SER:CB	1:B:146:PRO:HA	2.36	0.53
1:B:116:LEU:HD23	1:B:118:ILE:CG1	2.38	0.53
1:B:300:PRO:HB2	1:B:301:PRO:HD2	1.91	0.53
1:A:180:MET:CE	1:A:227:LEU:HD21	2.39	0.52
1:A:252:ASN:HD21	1:A:271:GLU:H	1.57	0.52
1:A:229:HIS:HD2	1:A:252:ASN:H	1.56	0.52
1:B:207:HIS:H	1:B:225:GLU:HG2	1.74	0.52
1:B:176:VAL:O	1:B:180:MET:HG2	2.09	0.52
1:B:226:PHE:CE1	1:B:252:ASN:HB3	2.45	0.51
1:B:201:LEU:HD21	1:B:204:LYS:HB3	1.93	0.51
1:A:138:LEU:HD13	1:A:192:TYR:CE2	2.46	0.50
1:B:40:SER:HB3	1:B:59:VAL:HB	1.93	0.50
1:B:150:LEU:HD23	1:B:166:ASP:HB2	1.93	0.50
1:A:141:SER:CB	1:A:146:PRO:HA	2.40	0.50
1:B:56:TYR:CE1	1:B:66:LYS:HD3	2.47	0.50
1:B:187:GLY:HA3	1:B:209:PRO:CD	2.41	0.50
1:A:226:PHE:CE1	2:A:1334:S55:H101	2.46	0.49
1:B:317:LEU:HD13	1:B:330:ILE:HG21	1.93	0.49
1:A:126:SER:HB2	1:A:136:THR:O	2.13	0.49
1:B:286:LYS:HG2	1:B:294:LEU:HB2	1.95	0.48
1:A:116:LEU:HD12	1:A:116:LEU:C	2.30	0.48
1:A:229:HIS:HD2	1:A:251:PRO:HA	1.78	0.48
1:B:235:TRP:HZ3	1:B:245:GLU:CD	2.17	0.48
1:A:241:LYS:HE3	3:A:2030:HOH:O	2.14	0.48
1:B:122:TYR:OH	1:B:173:ASP:HB2	2.14	0.48
1:B:220:PHE:CE1	1:B:222:LEU:HD21	2.48	0.48
1:A:40:SER:HB3	1:A:59:VAL:HB	1.94	0.47
1:A:211:GLY:CA	1:A:255:ASN:HA	2.43	0.47
1:B:209:PRO:HA	1:B:224:ALA:O	2.14	0.47
1:A:221:VAL:CG2	1:A:236:LEU:HD11	2.41	0.47
1:A:94:ASP:N	1:A:94:ASP:OD1	2.46	0.47
2:B:1334:S55:H9	2:B:1334:S55:H15	1.79	0.47
1:B:266:VAL:HG22	1:B:267:SER:N	2.29	0.47
1:A:96:GLU:OE1	1:A:96:GLU:N	2.39	0.47
1:A:98:ARG:N	1:A:99:PRO:CD	2.79	0.46
1:B:149:TRP:CD1	1:B:150:LEU:O	2.68	0.46
1:B:332:VAL:O	1:B:332:VAL:HG12	2.15	0.46
1:A:189:LEU:HD13	1:A:223:VAL:HB	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:SER:HB2	1:B:123:TYR:CE2	2.51	0.46
1:A:67:TYR:CE2	1:A:69:GLY:HA2	2.51	0.46
1:B:226:PHE:HA	1:B:253:PRO:HD2	1.98	0.45
1:A:108:SER:OG	1:A:155:VAL:HG23	2.17	0.45
1:B:229:HIS:CD2	1:B:252:ASN:N	2.70	0.45
1:B:207:HIS:HD2	1:B:228:SER:OG	2.00	0.45
1:B:253:PRO:HA	1:B:267:SER:O	2.16	0.45
1:A:180:MET:HE1	1:A:227:LEU:HD21	1.98	0.45
1:A:46:PHE:HA	1:A:56:TYR:O	2.18	0.44
1:B:56:TYR:CD1	1:B:66:LYS:HB2	2.53	0.44
1:B:113:ASN:ND2	1:B:115:GLN:HG3	2.33	0.44
1:B:44:ASN:OD1	1:B:58:SER:OG	2.36	0.43
1:B:71:ASN:ND2	1:B:71:ASN:H	2.15	0.43
1:A:265:TRP:CD2	1:A:286:LYS:HB2	2.53	0.43
1:B:233:LYS:HD3	1:B:235:TRP:CZ2	2.53	0.43
1:A:138:LEU:HD23	1:A:138:LEU:HA	1.81	0.43
1:A:180:MET:HE1	1:A:226:PHE:HE2	1.83	0.43
1:B:163:TYR:CD2	1:B:163:TYR:N	2.86	0.43
1:A:49:ASP:OD1	1:A:49:ASP:C	2.57	0.43
1:B:46:PHE:HA	1:B:56:TYR:O	2.19	0.43
1:B:121:CYS:HB3	1:B:122:TYR:CD2	2.54	0.43
1:B:275:ASN:HD22	1:B:277:HIS:H	1.66	0.42
1:A:95:ALA:HB1	1:A:174:ARG:HD2	2.01	0.42
1:A:230:GLN:HE21	1:A:230:GLN:HB3	1.66	0.42
1:A:214:VAL:HG13	1:A:215:SER:O	2.19	0.42
1:A:322:THR:HG21	1:A:325:HIS:HB2	2.01	0.42
1:B:88:PHE:CE1	1:B:97:LYS:HD3	2.53	0.42
1:B:281:ASP:HB3	1:B:283:LYS:NZ	2.33	0.42
1:B:61:ASP:O	1:B:84:TRP:NE1	2.45	0.42
1:B:210:GLY:CA	2:B:1334:S55:H1	2.49	0.42
1:A:89:CYS:O	1:A:92:SER:HB3	2.20	0.42
1:B:229:HIS:HD2	1:B:251:PRO:CA	2.25	0.42
1:A:113:ASN:HD22	1:A:113:ASN:C	2.23	0.42
1:A:84:TRP:O	1:A:85:ASN:ND2	2.53	0.41
1:A:235:TRP:HZ3	1:A:245:GLU:OE2	2.03	0.41
1:B:235:TRP:HZ3	1:B:245:GLU:HG2	1.84	0.41
1:B:233:LYS:HD3	1:B:235:TRP:CH2	2.55	0.41
1:A:142:VAL:CG2	1:A:199:THR:HG22	2.48	0.41
1:B:266:VAL:CG2	1:B:267:SER:N	2.84	0.41
1:B:98:ARG:HD3	1:B:173:ASP:OD1	2.20	0.41
1:B:279:ARG:HE	1:B:279:ARG:HB2	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:PRO:CB	1:B:301:PRO:HD2	2.50	0.41
1:A:189:LEU:HB3	1:A:203:LEU:HB2	2.01	0.41
2:B:1334:S55:H15	2:B:1334:S55:N9	2.36	0.41
1:A:215:SER:HB3	1:A:220:PHE:CZ	2.56	0.40
1:B:113:ASN:ND2	1:B:115:GLN:CG	2.84	0.40
1:A:253:PRO:HA	1:A:267:SER:O	2.22	0.40
1:A:251:PRO:O	1:A:252:ASN:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	300/302 (99%)	281 (94%)	16 (5%)	3 (1%)	15 50
1	B	300/302 (99%)	282 (94%)	17 (6%)	1 (0%)	41 75
All	All	600/604 (99%)	563 (94%)	33 (6%)	4 (1%)	22 59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	173	ASP
1	A	173	ASP
1	A	71	ASN
1	A	301	PRO

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	243/260 (94%)	213 (88%)	30 (12%)	4 20
1	B	243/260 (94%)	210 (86%)	33 (14%)	3 16
All	All	486/520 (94%)	423 (87%)	63 (13%)	4 18

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LYS
1	A	47	THR
1	A	53	LYS
1	A	68	GLU
1	A	72	SER
1	A	85	ASN
1	A	86	LYS
1	A	92	SER
1	A	96	GLU
1	A	103	ARG
1	A	113	ASN
1	A	131	GLU
1	A	141	SER
1	A	142	VAL
1	A	155	VAL
1	A	158	ARG
1	A	165	THR
1	A	170	LEU
1	A	173	ASP
1	A	174	ARG
1	A	176	VAL
1	A	178	GLN
1	A	191	LYS
1	A	203	LEU
1	A	207	HIS
1	A	214	VAL
1	A	221	VAL
1	A	235	TRP
1	A	301	PRO
1	A	308	PHE
1	B	35	LEU
1	B	51	THR
1	B	66	LYS

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Mol	Chain	Res	Type
1	B	71	ASN
1	B	75	VAL
1	B	92	SER
1	B	100	LEU
1	B	113	ASN
1	B	116	LEU
1	B	126	SER
1	B	128	VAL
1	B	142	VAL
1	B	148	LYS
1	B	158	ARG
1	B	165	THR
1	B	173	ASP
1	B	174	ARG
1	B	181	ASP
1	B	183	SER
1	B	186	THR
1	B	195	SER
1	B	197	LYS
1	B	199	THR
1	B	200	THR
1	B	207	HIS
1	B	219	SER
1	B	221	VAL
1	B	228	SER
1	B	235	TRP
1	B	248	VAL
1	B	275	ASN
1	B	283	LYS
1	B	333	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	85	ASN
1	A	91	ASN
1	A	113	ASN
1	A	178	GLN
1	A	207	HIS
1	A	229	HIS
1	A	230	GLN

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Mol	Chain	Res	Type
1	A	252	ASN
1	A	275	ASN
1	A	277	HIS
1	B	91	ASN
1	B	113	ASN
1	B	114	ASN
1	B	207	HIS
1	B	229	HIS
1	B	230	GLN
1	B	252	ASN
1	B	275	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

2 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	$\#\|Z\| > 2$
2	S55	B	1334	-	39,42,42	1.88	6 (15%)	44,61,61	2.25	12 (27%)
2	S55	A	1334	-	39,42,42	1.66	3 (7%)	44,61,61	2.72	16 (36%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	S55	B	1334	-	-	1/18/65/65	0/5/5/5
2	S55	A	1334	-	-	3/18/65/65	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1334	S55	O18-C19	6.78	1.48	1.35
2	A	1334	S55	O18-C19	6.09	1.46	1.35
2	B	1334	S55	O22-C21	6.08	1.46	1.33
2	A	1334	S55	O22-C21	5.75	1.46	1.33
2	B	1334	S55	C14-C13	-3.43	1.50	1.53
2	A	1334	S55	C13-N12	-3.40	1.44	1.47
2	B	1334	S55	C13-N12	-2.51	1.45	1.47
2	B	1334	S55	C2-C3	2.26	1.41	1.36
2	B	1334	S55	C17-C16	2.12	1.55	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1334	S55	O22-C21-C20	9.12	124.52	112.64
2	B	1334	S55	O18-C17-C16	-7.82	101.86	113.16
2	A	1334	S55	O18-C17-C16	-6.72	103.44	113.16
2	B	1334	S55	O22-C21-C20	6.41	120.98	112.64
2	A	1334	S55	C16-C15-C20	-5.07	104.30	108.60
2	A	1334	S55	O24-C21-C20	-4.53	118.67	124.34
2	A	1334	S55	C10-C11-N12	4.30	115.00	109.04
2	A	1334	S55	C23-O22-C21	4.24	123.87	115.86
2	A	1334	S55	C11-N12-C13	4.00	119.71	111.69
2	B	1334	S55	C11-N12-C13	3.87	119.46	111.69
2	B	1334	S55	C23-O22-C21	3.77	123.00	115.86
2	A	1334	S55	O22-C21-O24	-3.54	116.72	123.53
2	A	1334	S55	C17-O18-C19	3.39	122.69	116.40
2	B	1334	S55	C6-C5-C4	3.24	122.47	118.17
2	A	1334	S55	C6-C5-C4	3.05	122.22	118.17
2	B	1334	S55	C17-O18-C19	2.99	121.95	116.40
2	B	1334	S55	O22-C21-O24	-2.95	117.86	123.53
2	A	1334	S55	O18-C19-C20	-2.87	120.90	125.58
2	A	1334	S55	C28-O33-C32	2.62	118.84	113.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1334	S55	O36-C30-C31	-2.54	104.47	110.35
2	B	1334	S55	O24-C21-C20	-2.53	121.16	124.34
2	A	1334	S55	C6-C5-C7	-2.47	129.92	134.17
2	A	1334	S55	O33-C32-C33	2.37	112.32	106.44
2	B	1334	S55	O27-C28-O33	-2.26	104.37	110.67
2	A	1334	S55	O35-C31-C32	-2.22	103.77	109.30
2	B	1334	S55	O18-C19-C20	-2.20	122.00	125.58
2	B	1334	S55	C6-C5-C7	-2.15	130.46	134.17
2	A	1334	S55	O35-C31-C30	-2.08	105.54	110.35

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1334	S55	O24-C21-O22-C23
2	A	1334	S55	C20-C21-O22-C23
2	B	1334	S55	C15-C16-C25-C26
2	A	1334	S55	C13-C14-C15-C20

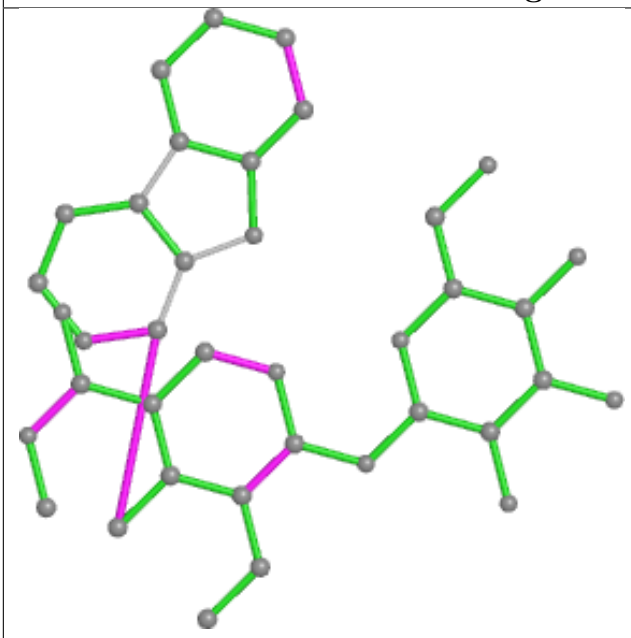
There are no ring outliers.

2 monomers are involved in 11 short contacts:

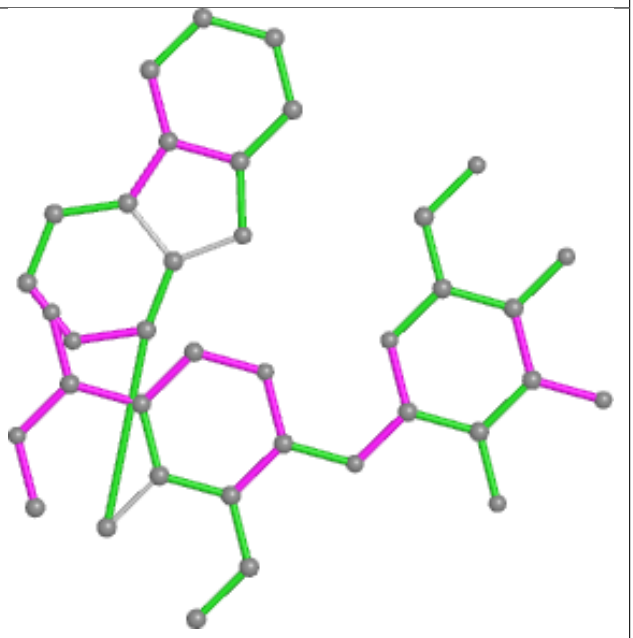
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1334	S55	4	0
2	A	1334	S55	7	0

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 then 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.

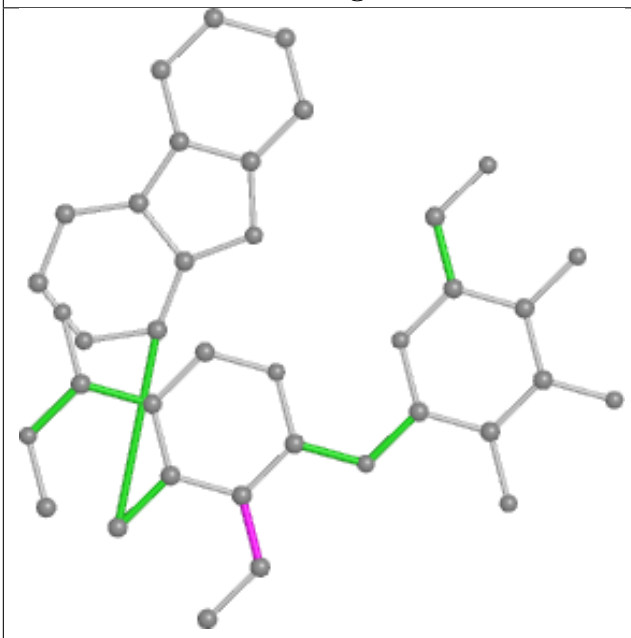
Ligand S55 B 1334



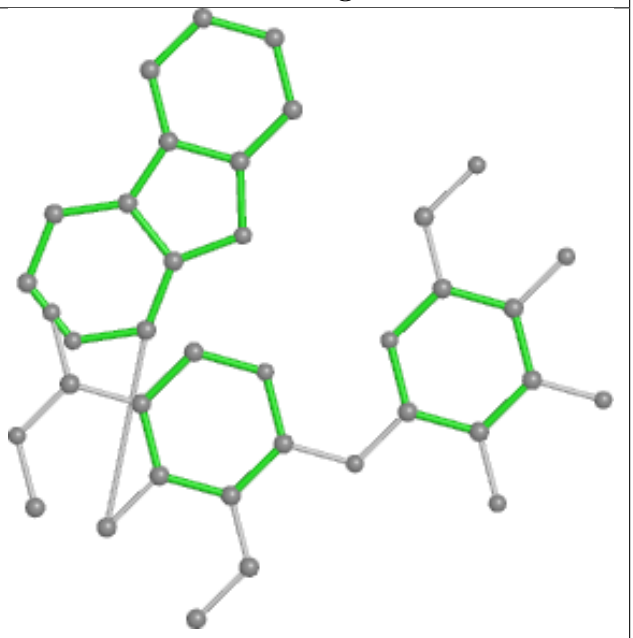
Bond lengths



Bond angles

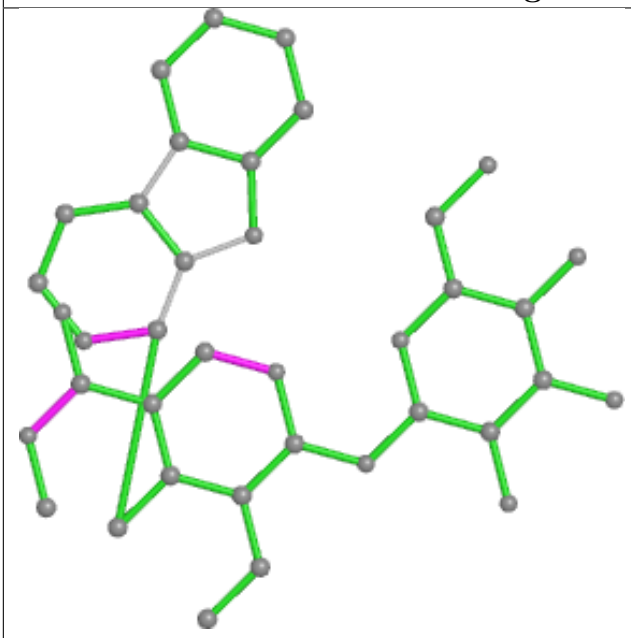


Torsions

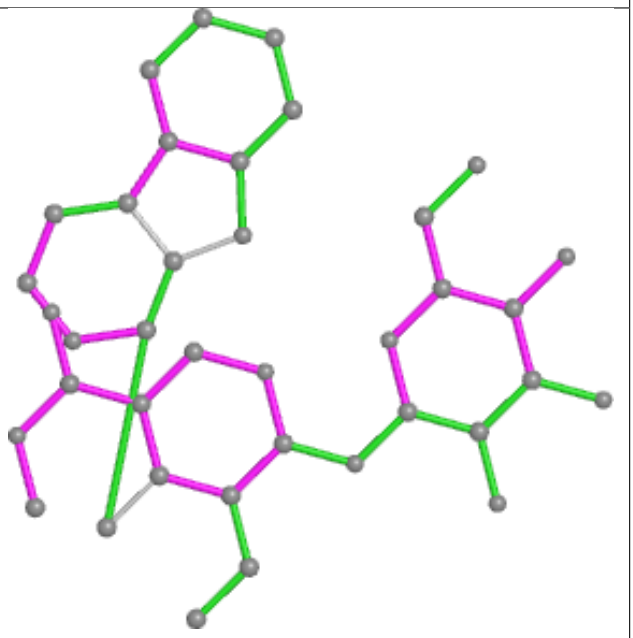


Rings

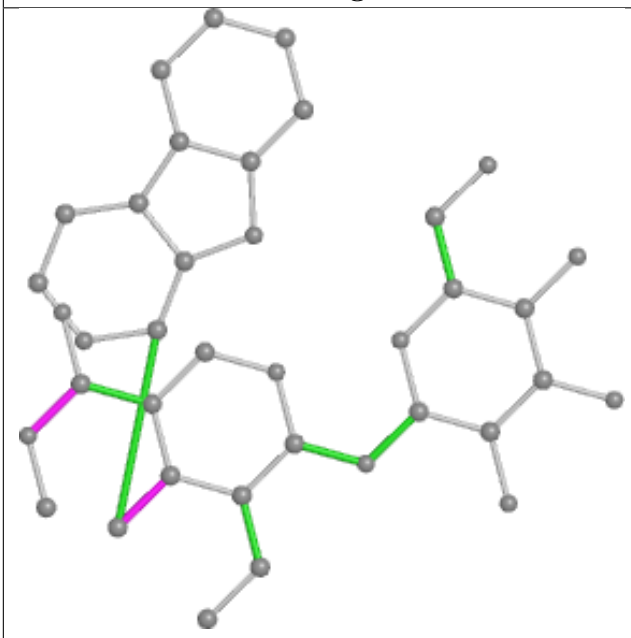
Ligand S55 A 1334



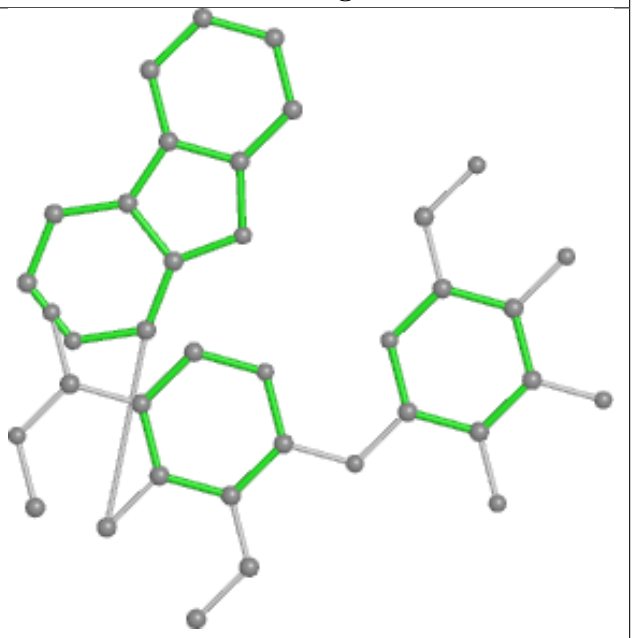
Bond lengths



Bond angles



Torsions



Rings

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^{th} 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	302/302 (100%)	-0.71	0	100	100	21, 26, 31, 42	0
1	B	302/302 (100%)	-0.58	0	100	100	20, 26, 31, 35	0
All	All	604/604 (100%)	-0.64	0	100	100	20, 26, 31, 42	0

There are no RSRZ outliers to report.

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

There are no monosaccharides in this entry.

6.4 Ligands

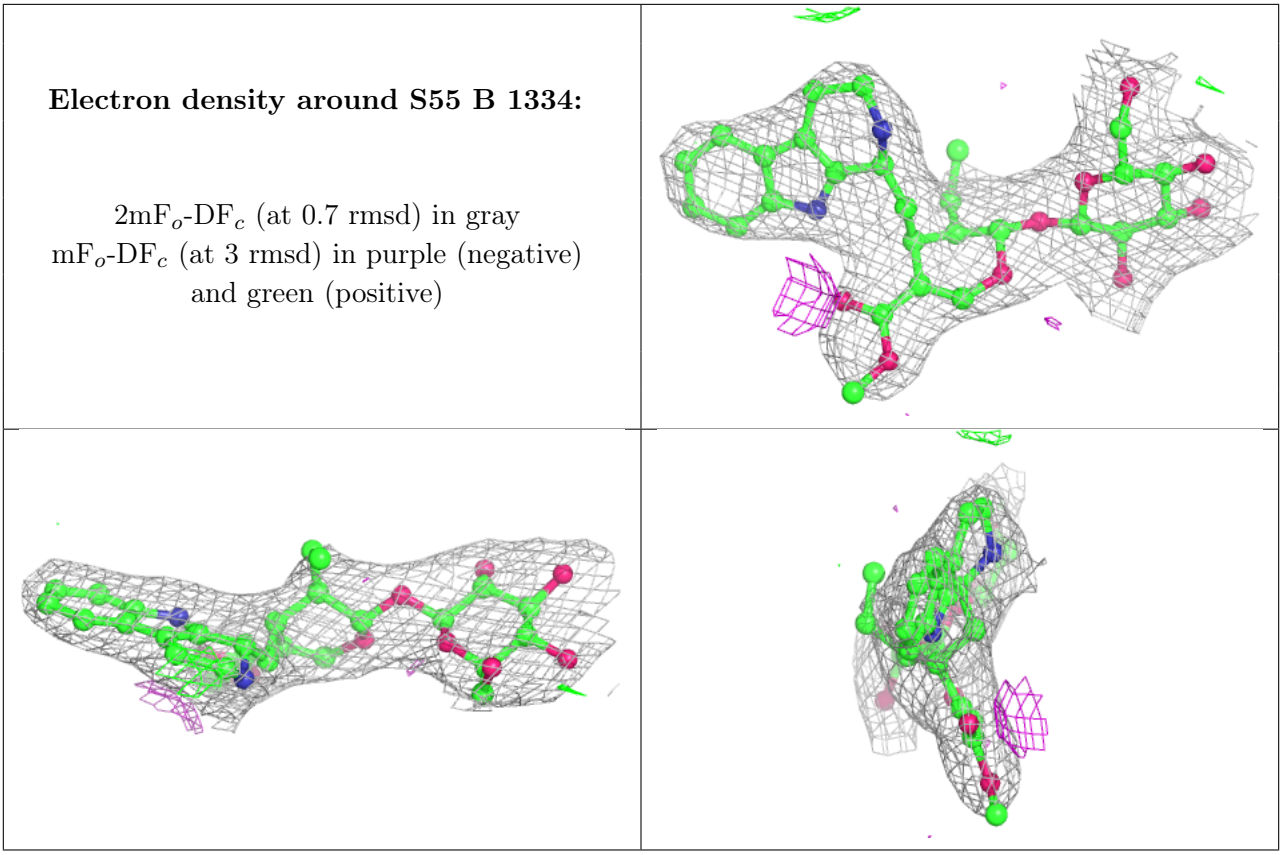
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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	S55	B	1334	38/38	0.95	0.14	45,53,62,64	0
2	S55	A	1334	38/38	0.97	0.13	42,48,56,58	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.

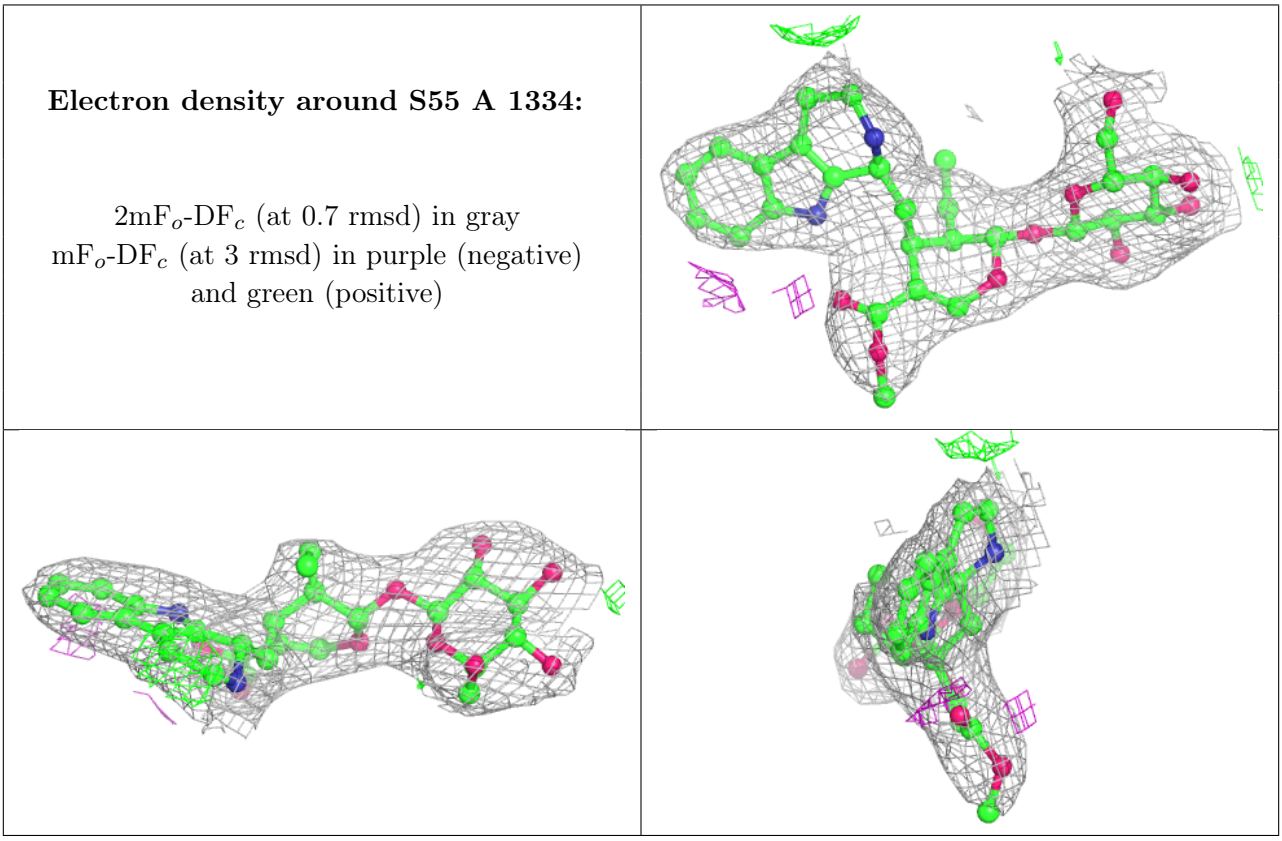
Electron density around S55 B 1334:

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



Electron density around S55 A 1334:

$2mF_o$-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.