



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:14 PM UTC

PDB ID : 8W71 / pdb_00008w71
Title : Structural basis of chorismate isomerization by Arabidopsis isochorismate synthase ICS1
Authors : Su, Z.H.; Ming, Z.H.
Deposited on : 2023-08-30
Resolution : 2.12 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

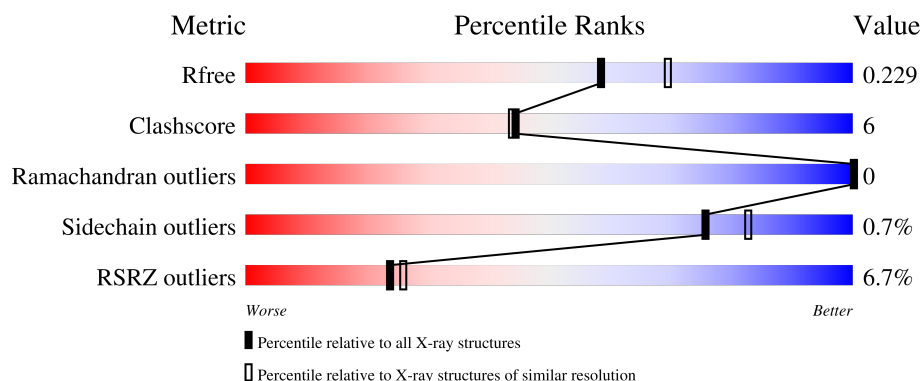
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.12 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	536	7% 81% 11% 8%
1	B	536	6% 80% 11% 8%

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	FMT	A	603	-	X	X	-
4	FMT	B	603	-	X	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8469 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 Isochorismate synthase 1, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	494	Total	C	N	O	S	0	0	0
			3836	2423	671	727	15			
1	B	493	Total	C	N	O	S	0	0	0
			3825	2417	667	726	15			

There are 26 discrepancies between the modelled and reference sequences:

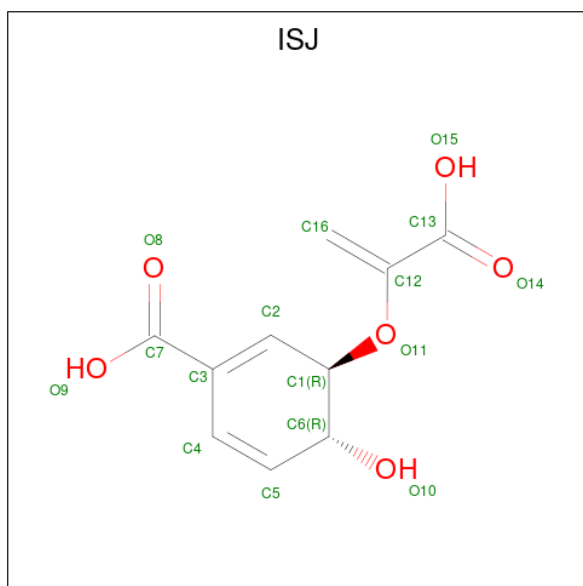
Chain	Residue	Modelled	Actual	Comment	Reference
A	44	MET	-	initiating methionine	UNP Q9S7H8
A	45	ASP	-	expression tag	UNP Q9S7H8
A	316	ARG	LYS	engineered mutation	UNP Q9S7H8
A	570	HIS	-	expression tag	UNP Q9S7H8
A	571	HIS	-	expression tag	UNP Q9S7H8
A	572	HIS	-	expression tag	UNP Q9S7H8
A	573	HIS	-	expression tag	UNP Q9S7H8
A	574	HIS	-	expression tag	UNP Q9S7H8
A	575	HIS	-	expression tag	UNP Q9S7H8
A	576	HIS	-	expression tag	UNP Q9S7H8
A	577	HIS	-	expression tag	UNP Q9S7H8
A	578	HIS	-	expression tag	UNP Q9S7H8
A	579	HIS	-	expression tag	UNP Q9S7H8
B	44	MET	-	initiating methionine	UNP Q9S7H8
B	45	ASP	-	expression tag	UNP Q9S7H8
B	316	ARG	LYS	engineered mutation	UNP Q9S7H8
B	570	HIS	-	expression tag	UNP Q9S7H8
B	571	HIS	-	expression tag	UNP Q9S7H8
B	572	HIS	-	expression tag	UNP Q9S7H8
B	573	HIS	-	expression tag	UNP Q9S7H8
B	574	HIS	-	expression tag	UNP Q9S7H8
B	575	HIS	-	expression tag	UNP Q9S7H8
B	576	HIS	-	expression tag	UNP Q9S7H8
B	577	HIS	-	expression tag	UNP Q9S7H8
B	578	HIS	-	expression tag	UNP Q9S7H8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	579	HIS	-	expression tag	UNP Q9S7H8

- Molecule 2 is (3R,4R)-3-[(1-carboxyethenyl)oxy]-4-hydroxycyclohexa-1,5-diene-1-carboxylic acid (CCD ID: ISJ) (formula: C₁₀H₁₀O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	1
			16	10	6		
2	B	1	Total	C	O	0	1
			16	10	6		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	B	1	Total	C	O	0	0
			3	1	2		

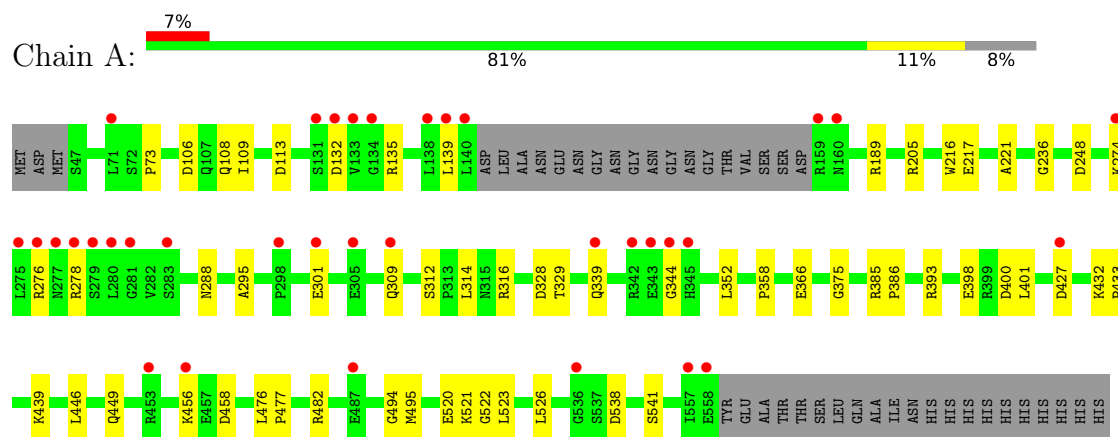
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	385	Total	O	0	0
			385	385		
5	B	383	Total	O	0	0
			383	383		

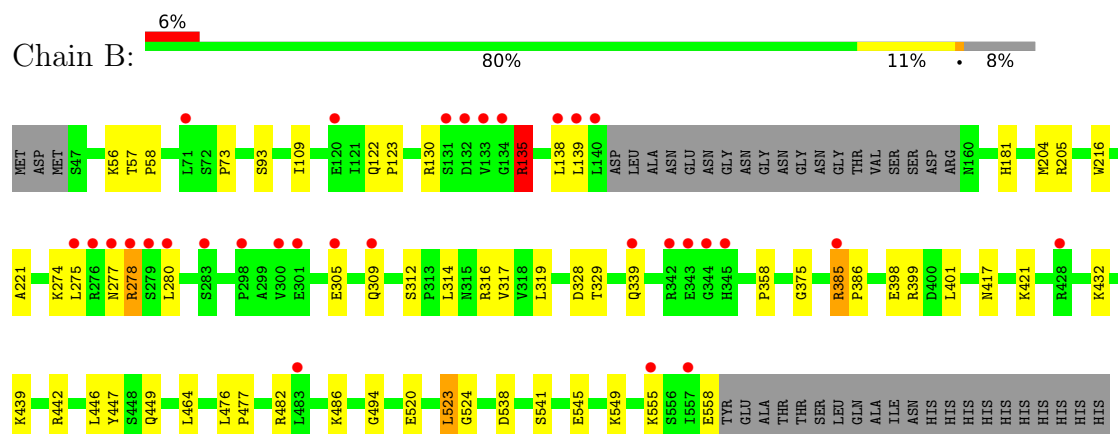
3 Residue-property plots [i](#)

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: Isochorismate synthase 1, chloroplastic



- Molecule 1: Isochorismate synthase 1, chloroplastic



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.10Å 73.90Å 103.78Å 70.84° 84.46° 90.02°	Depositor
Resolution (Å)	39.03 – 2.12 39.03 – 2.12	Depositor EDS
% Data completeness (in resolution range)	93.3 (39.03-2.12) 93.3 (39.03-2.12)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.12Å)	Xtriage
Refinement program	PHENIX 1.15_3459	Depositor
R, R_{free}	0.180 , 0.210 (Not available) , 0.229	Depositor DCC
R_{free} test set	4228 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.900	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.127 for -h,k,k-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8469	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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: FMT, MG, ISJ

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	0.48	4/3913 (0.1%)	0.62	4/5302 (0.1%)
1	B	0.52	4/3902 (0.1%)	0.67	5/5288 (0.1%)
All	All	0.50	8/7815 (0.1%)	0.65	9/10590 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	385	ARG	C-N	11.99	1.48	1.33
1	B	135	ARG	C-N	11.35	1.48	1.33
1	B	524	GLY	C-N	-7.69	1.24	1.33
1	A	526	LEU	C-O	-6.23	1.16	1.24
1	A	522	GLY	C-O	-5.51	1.16	1.24
1	B	386	PRO	C-O	-5.34	1.17	1.23
1	B	130	ARG	C-O	-5.28	1.17	1.23
1	A	386	PRO	C-O	-5.06	1.17	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	135	ARG	CA-C-N	10.72	130.82	119.78
1	B	135	ARG	C-N-CA	10.72	130.82	119.78
1	A	385	ARG	O-C-N	-10.65	111.24	121.57
1	A	385	ARG	CA-C-N	7.73	127.68	120.03
1	A	385	ARG	C-N-CA	7.73	127.68	120.03
1	B	278	ARG	N-CA-C	-6.19	105.86	113.41
1	B	135	ARG	O-C-N	-5.88	114.85	121.43
1	A	385	ARG	N-CA-C	-5.75	101.29	110.10
1	B	385	ARG	CB-CG-CD	5.34	123.59	111.30

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	3836	0	3827	45	0
1	B	3825	0	3814	48	0
2	A	16	0	8	0	0
2	B	16	0	8	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	2	2	0
4	B	3	0	2	2	0
5	A	385	0	0	16	2
5	B	383	0	0	8	2
All	All	8469	0	7661	92	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (92) 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:B:385:ARG:NH1	5:B:701:HOH:O	1.77	1.14
1:B:442:ARG:HH12	1:B:555:LYS:HD3	0.99	1.05
1:A:456:LYS:HE2	1:A:458:ASP:HB2	1.47	0.93
1:B:538:ASP:OD1	5:B:702:HOH:O	1.86	0.91
1:B:442:ARG:NH1	1:B:555:LYS:HD3	1.84	0.90
1:B:56:LYS:NZ	5:B:705:HOH:O	2.07	0.88
1:A:295:ALA:O	5:A:701:HOH:O	1.98	0.82
1:B:482:ARG:O	5:B:703:HOH:O	1.97	0.82
1:A:217:GLU:OE2	5:A:702:HOH:O	1.98	0.81
1:B:442:ARG:HH22	1:B:555:LYS:HE3	1.45	0.79
1:A:312:SER:O	5:A:704:HOH:O	2.05	0.74
1:A:352:LEU:HD23	1:A:495:MET:HE1	1.70	0.73
1:A:482:ARG:O	5:A:705:HOH:O	2.06	0.72
1:B:316:ARG:HH12	4:B:603:FMT:C	2.04	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:ASP:OD2	5:A:707:HOH:O	2.11	0.69
1:A:427:ASP:O	5:A:706:HOH:O	2.10	0.69
1:A:366:GLU:OE2	5:A:709:HOH:O	2.14	0.66
1:A:400:ASP:OD1	5:A:708:HOH:O	2.14	0.66
1:A:456:LYS:CE	1:A:458:ASP:HB2	2.26	0.63
1:A:482:ARG:NH1	5:A:703:HOH:O	2.03	0.62
1:B:538:ASP:OD2	1:B:541:SER:HB2	2.00	0.62
1:A:276:ARG:HH11	1:A:276:ARG:HG2	1.65	0.62
1:A:521:LYS:O	5:A:710:HOH:O	2.16	0.61
1:B:181:HIS:HD2	5:B:720:HOH:O	1.84	0.61
1:B:442:ARG:HH12	1:B:555:LYS:CD	1.93	0.60
1:A:288:ASN:ND2	5:A:714:HOH:O	2.25	0.60
1:B:278:ARG:HH11	1:B:278:ARG:HG3	1.66	0.60
1:A:108:GLN:HG2	1:A:236:GLY:C	2.29	0.58
1:B:482:ARG:HH12	1:B:486:LYS:HE2	1.69	0.57
1:B:277:ASN:O	1:B:339:GLN:OE1	2.25	0.55
1:B:305:GLU:HB3	1:B:309:GLN:OE1	2.07	0.54
1:B:442:ARG:NH2	1:B:555:LYS:HE3	2.20	0.54
1:B:205:ARG:HD3	1:B:221:ALA:HA	1.90	0.54
1:A:109:ILE:HG23	1:A:274:LYS:HZ1	1.73	0.54
1:B:482:ARG:NH1	1:B:486:LYS:HE2	2.23	0.53
1:A:106:ASP:O	5:A:711:HOH:O	2.19	0.52
1:A:276:ARG:HG2	1:A:276:ARG:NH1	2.23	0.52
1:A:309:GLN:O	5:A:704:HOH:O	2.19	0.52
1:A:316:ARG:HH12	4:A:603:FMT:C	2.24	0.51
1:B:316:ARG:HH12	4:B:603:FMT:H	1.75	0.51
1:A:316:ARG:HH12	4:A:603:FMT:H	1.76	0.51
1:B:278:ARG:HG3	1:B:278:ARG:NH1	2.26	0.50
1:A:358:PRO:HB3	1:A:520:GLU:HG2	1.92	0.50
1:B:216:TRP:CD2	1:B:494:GLY:HA3	2.47	0.49
1:A:398:GLU:OE1	1:A:439:LYS:HE2	2.13	0.49
1:B:73:PRO:HG3	1:B:109:ILE:HD13	1.95	0.49
1:B:401:LEU:HD22	1:B:446:LEU:HG	1.95	0.48
1:A:278:ARG:HD2	1:A:339:GLN:NE2	2.28	0.48
1:A:344:GLY:O	5:A:712:HOH:O	2.20	0.48
1:B:358:PRO:HB3	1:B:520:GLU:HG2	1.94	0.48
1:B:482:ARG:HH12	1:B:486:LYS:CE	2.27	0.48
1:A:139:LEU:HD22	1:A:375:GLY:HA3	1.96	0.47
1:B:93:SER:OG	5:B:706:HOH:O	2.19	0.47
1:A:132:ASP:OD1	1:A:135:ARG:NH2	2.48	0.46
1:A:538:ASP:OD2	1:A:541:SER:HB2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:VAL:HG13	1:B:319:LEU:HD21	1.98	0.46
1:A:109:ILE:HG23	1:A:274:LYS:NZ	2.31	0.46
1:B:312:SER:O	5:B:707:HOH:O	2.20	0.46
1:B:216:TRP:CE2	1:B:494:GLY:HA3	2.51	0.45
1:B:204:MET:HE1	1:B:464:LEU:HD21	1.97	0.45
1:B:314:LEU:HA	1:B:477:PRO:HD2	1.97	0.45
1:A:216:TRP:CD2	1:A:494:GLY:HA3	2.53	0.44
1:B:139:LEU:HD22	1:B:375:GLY:HA3	1.99	0.44
1:A:401:LEU:HG	1:A:446:LEU:HG	2.00	0.43
1:B:122:GLN:HA	1:B:123:PRO:C	2.42	0.43
1:B:417:ASN:O	1:B:421:LYS:HG2	2.18	0.43
1:B:317:VAL:HG13	1:B:319:LEU:CD2	2.49	0.43
1:B:545:GLU:HG2	1:B:549:LYS:HE2	2.01	0.43
1:A:432:LYS:HA	1:A:433:PRO:HA	1.91	0.43
1:B:309:GLN:HB2	5:B:774:HOH:O	2.17	0.43
1:A:393:ARG:HD2	1:A:393:ARG:HA	1.82	0.43
1:A:476:LEU:HA	1:A:477:PRO:C	2.44	0.43
1:B:204:MET:HE1	1:B:464:LEU:HD11	2.01	0.42
1:B:138:LEU:HD12	1:B:138:LEU:HA	1.76	0.42
1:B:328:ASP:OD1	1:B:329:THR:N	2.45	0.42
1:B:555:LYS:O	1:B:558:GLU:HG3	2.20	0.42
1:A:314:LEU:HA	1:A:477:PRO:HD2	2.02	0.42
1:A:205:ARG:HD3	1:A:221:ALA:HA	2.00	0.42
1:A:113:ASP:OD2	1:A:274:LYS:HD2	2.20	0.41
1:B:398:GLU:OE1	1:B:439:LYS:HE2	2.19	0.41
1:B:432:LYS:O	1:B:449:GLN:HB2	2.20	0.41
1:A:328:ASP:O	1:B:399:ARG:NH2	2.53	0.41
1:A:432:LYS:O	1:A:449:GLN:HB2	2.20	0.41
1:A:328:ASP:OD1	1:A:329:THR:N	2.52	0.41
1:A:73:PRO:HG3	1:A:109:ILE:HD13	2.01	0.41
1:B:523:LEU:HD23	1:B:523:LEU:HA	1.74	0.41
1:B:476:LEU:HA	1:B:477:PRO:C	2.46	0.41
1:A:301:GLU:OE2	5:A:713:HOH:O	2.20	0.40
1:B:57:THR:HA	1:B:58:PRO:HD3	1.96	0.40
1:A:189:ARG:HD3	5:A:822:HOH:O	2.19	0.40
1:B:135:ARG:HD2	1:B:447:TYR:HE2	1.87	0.40
1:A:216:TRP:CE2	1:A:494:GLY:HA3	2.56	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:970:HOH:O	5:B:994:HOH:O[1_655]	1.94	0.26
5:A:730:HOH:O	5:A:946:HOH:O[1_655]	2.08	0.12
5:A:1037:HOH:O	5:B:815:HOH:O[1_545]	2.16	0.04

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	490/536 (91%)	476 (97%)	14 (3%)	0	100	100
1	B	489/536 (91%)	475 (97%)	14 (3%)	0	100	100
All	All	979/1072 (91%)	951 (97%)	28 (3%)	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	415/450 (92%)	414 (100%)	1 (0%)	87	92
1	B	414/450 (92%)	409 (99%)	5 (1%)	63	71
All	All	829/900 (92%)	823 (99%)	6 (1%)	76	83

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	523	LEU

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Mol	Chain	Res	Type
1	B	135	ARG
1	B	274	LYS
1	B	275	LEU
1	B	280	LEU
1	B	523	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	288	ASN
1	A	309	GLN
1	A	417	ASN
1	A	449	GLN
1	B	160	ASN
1	B	181	HIS
1	B	417	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 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	$\# Z > 2$
2	ISJ	A	601[B]	-	16,16,16	4.40	8 (50%)	13,22,22	3.74	7 (53%)
4	FMT	A	603	-	2,2,2	3.02	2 (100%)	1,1,1	1.42	0
4	FMT	B	603	-	2,2,2	3.05	2 (100%)	1,1,1	1.31	0
2	ISJ	B	601[B]	-	16,16,16	4.39	8 (50%)	13,22,22	4.01	8 (61%)

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	ISJ	A	601[B]	-	-	6/10/25/25	0/1/1/1
2	ISJ	B	601[B]	-	-	5/10/25/25	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601[B]	ISJ	C4-C5	10.69	1.53	1.33
2	A	601[B]	ISJ	C4-C5	10.47	1.53	1.33
2	A	601[B]	ISJ	C2-C3	7.82	1.53	1.35
2	B	601[B]	ISJ	C2-C3	7.82	1.53	1.35
2	A	601[B]	ISJ	C16-C12	7.81	1.52	1.31
2	B	601[B]	ISJ	C16-C12	7.63	1.52	1.31
2	B	601[B]	ISJ	C4-C3	5.00	1.53	1.43
2	A	601[B]	ISJ	C4-C3	4.82	1.53	1.43
4	B	603	FMT	O1-C	3.57	1.41	1.22
4	A	603	FMT	O1-C	3.46	1.41	1.22
2	B	601[B]	ISJ	C1-C2	3.42	1.55	1.50
2	A	601[B]	ISJ	C3-C7	3.28	1.54	1.48
2	B	601[B]	ISJ	C3-C7	3.18	1.54	1.48
2	A	601[B]	ISJ	C12-C13	2.99	1.52	1.49
2	A	601[B]	ISJ	C1-C2	2.99	1.54	1.50
2	B	601[B]	ISJ	C12-C13	2.57	1.51	1.49
4	A	603	FMT	O2-C	2.49	1.41	1.28
4	B	603	FMT	O2-C	2.42	1.41	1.28
2	B	601[B]	ISJ	C6-C1	2.24	1.55	1.52
2	A	601[B]	ISJ	C6-C1	2.05	1.54	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601[B]	ISJ	C5-C4-C3	-9.09	110.18	122.38
2	A	601[B]	ISJ	C5-C4-C3	-8.32	111.22	122.38
2	A	601[B]	ISJ	C16-C12-C13	-7.57	108.73	122.73
2	B	601[B]	ISJ	C16-C12-C13	-6.89	110.00	122.73
2	B	601[B]	ISJ	C6-C5-C4	-4.68	113.81	123.08
2	A	601[B]	ISJ	C6-C5-C4	-4.64	113.89	123.08
2	B	601[B]	ISJ	O15-C13-C12	4.62	121.79	113.91
2	B	601[B]	ISJ	C1-O11-C12	-3.26	112.00	117.03
2	B	601[B]	ISJ	O14-C13-C12	-2.98	117.29	121.79
2	A	601[B]	ISJ	C1-O11-C12	-2.80	112.70	117.03
2	A	601[B]	ISJ	O9-C7-C3	2.55	120.47	113.72
2	A	601[B]	ISJ	O11-C12-C13	-2.30	109.89	115.59
2	A	601[B]	ISJ	O15-C13-C12	2.16	117.59	113.91
2	B	601[B]	ISJ	O11-C12-C13	-2.12	110.34	115.59
2	B	601[B]	ISJ	O9-C7-C3	2.11	119.32	113.72

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601[B]	ISJ	C2-C3-C7-O8
2	A	601[B]	ISJ	C2-C3-C7-O9
2	A	601[B]	ISJ	C16-C12-C13-O14
2	A	601[B]	ISJ	C16-C12-C13-O15
2	B	601[B]	ISJ	C2-C3-C7-O8
2	B	601[B]	ISJ	C2-C3-C7-O9
2	B	601[B]	ISJ	O11-C12-C13-O15
2	B	601[B]	ISJ	C16-C12-C13-O14
2	B	601[B]	ISJ	C16-C12-C13-O15
2	A	601[B]	ISJ	O11-C12-C13-O15
2	A	601[B]	ISJ	O11-C12-C13-O14

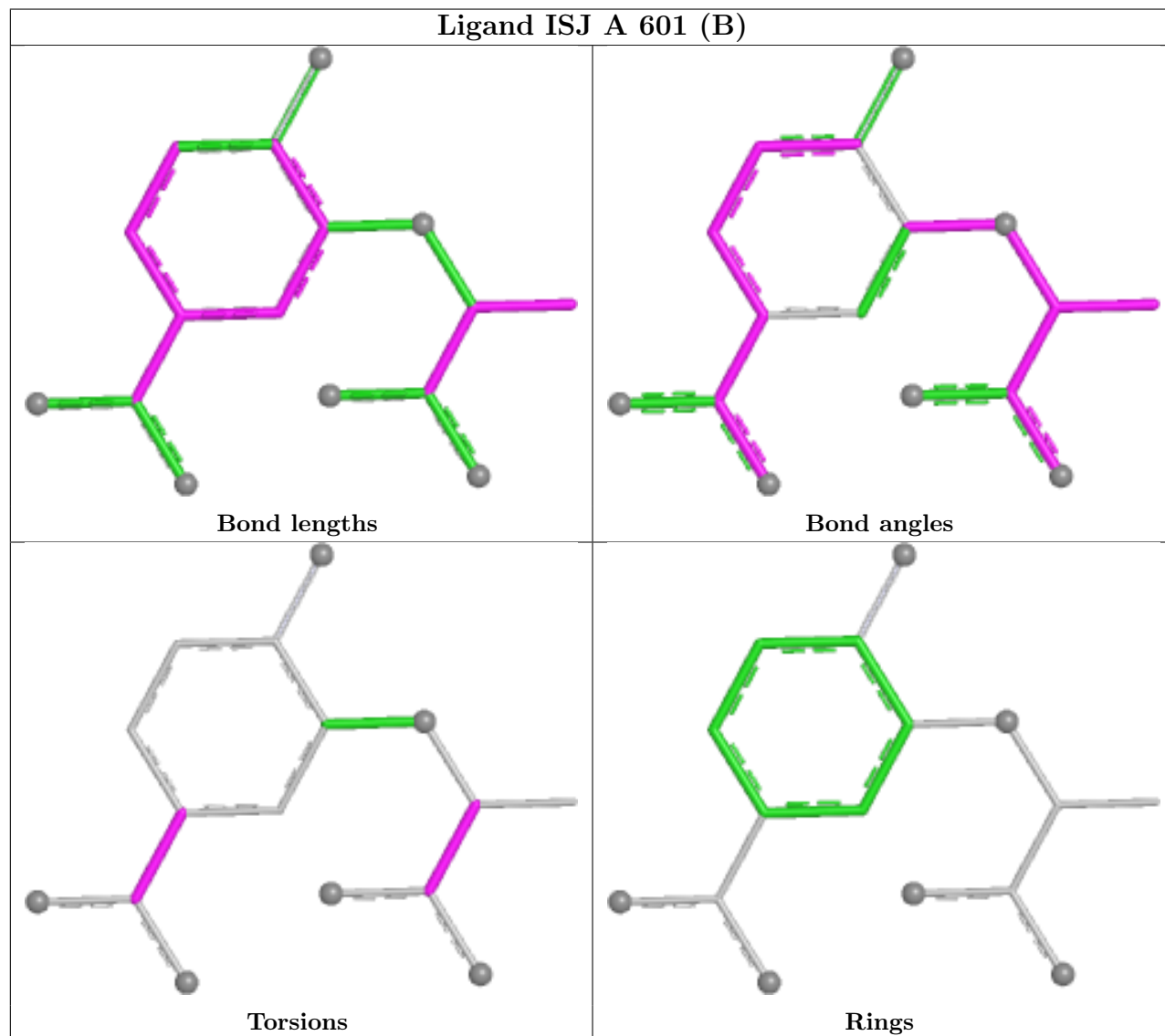
There are no ring outliers.

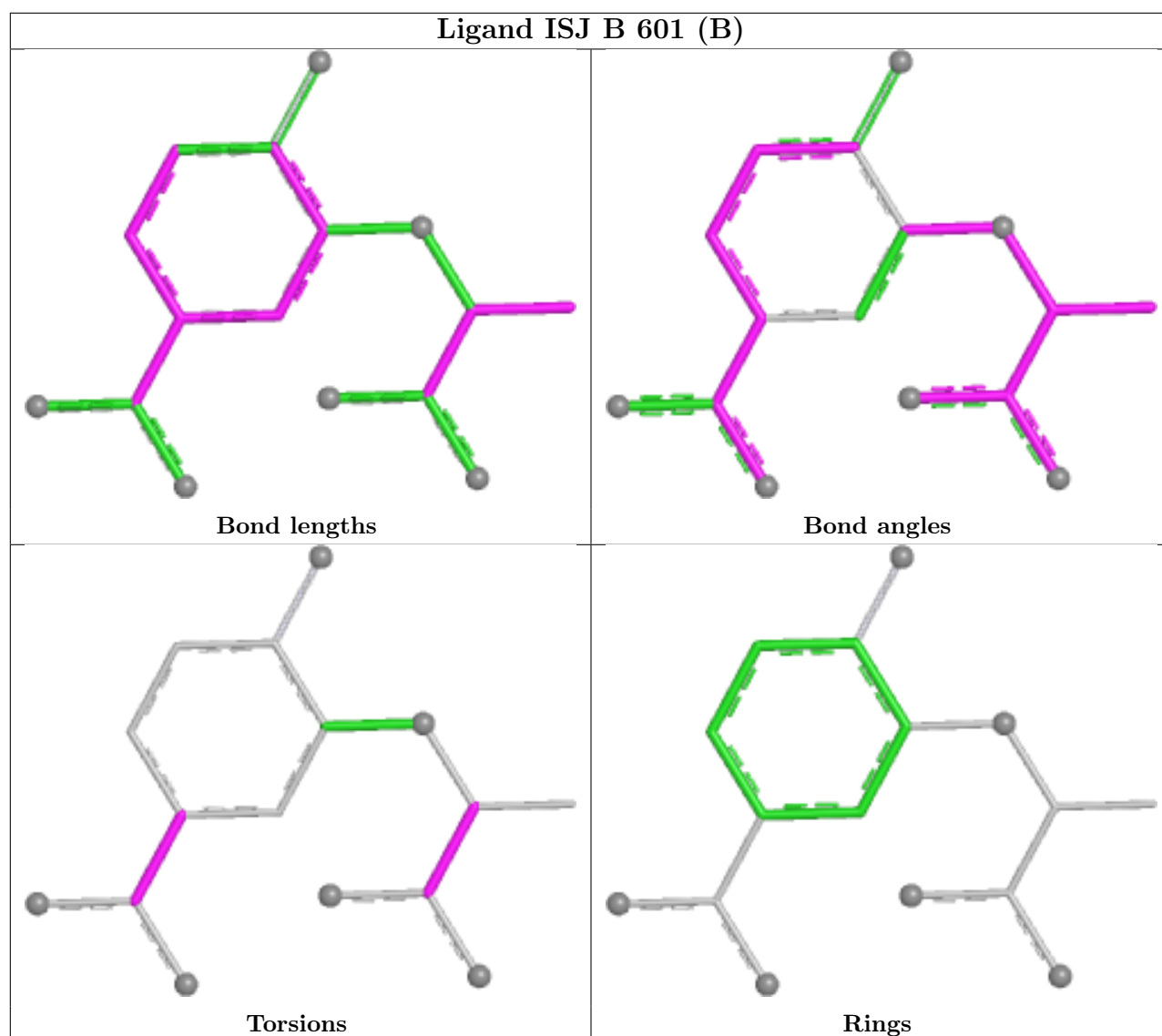
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	603	FMT	2	0
4	B	603	FMT	2	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	494/536 (92%)	0.20	35 (7%)	22 24	15, 26, 52, 82	0
1	B	493/536 (91%)	0.24	31 (6%)	26 28	16, 27, 52, 82	0
All	All	987/1072 (92%)	0.22	66 (6%)	24 26	15, 26, 52, 82	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	133	VAL	7.4
1	B	277	ASN	7.1
1	A	140	LEU	6.1
1	B	132	ASP	6.1
1	B	344	GLY	5.9
1	A	133	VAL	5.3
1	B	140	LEU	5.3
1	A	344	GLY	5.1
1	A	132	ASP	4.8
1	B	279	SER	4.6
1	A	345	HIS	4.6
1	A	275	LEU	4.6
1	B	278	ARG	4.4
1	A	557	ILE	4.4
1	B	557	ILE	4.1
1	B	275	LEU	4.1
1	B	305	GLU	4.0
1	A	279	SER	3.9
1	B	345	HIS	3.9
1	B	342	ARG	3.8
1	A	305	GLU	3.7
1	A	343	GLU	3.5
1	B	131	SER	3.4
1	A	276	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	139	LEU	3.4
1	A	280	LEU	3.1
1	A	427	ASP	3.1
1	B	339	GLN	3.1
1	A	277	ASN	3.0
1	A	453	ARG	3.0
1	B	120	GLU	3.0
1	A	278	ARG	2.9
1	B	283	SER	2.9
1	B	139	LEU	2.9
1	A	159	ARG	2.8
1	B	134	GLY	2.8
1	A	342	ARG	2.7
1	B	276	ARG	2.7
1	A	301	GLU	2.7
1	B	343	GLU	2.7
1	A	134	GLY	2.7
1	B	301	GLU	2.6
1	A	309	GLN	2.6
1	B	309	GLN	2.6
1	A	131	SER	2.6
1	A	283	SER	2.5
1	A	71	LEU	2.5
1	A	281	GLY	2.5
1	A	138	LEU	2.5
1	B	483	LEU	2.5
1	B	138	LEU	2.4
1	A	487	GLU	2.4
1	B	555	LYS	2.4
1	A	536	GLY	2.3
1	B	428	ARG	2.3
1	B	298	PRO	2.2
1	B	385	ARG	2.2
1	B	280	LEU	2.2
1	A	298	PRO	2.2
1	B	300	VAL	2.1
1	A	339	GLN	2.1
1	A	160	ASN	2.1
1	A	274	LYS	2.1
1	A	456	LYS	2.1
1	B	71	LEU	2.0
1	A	558	GLU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

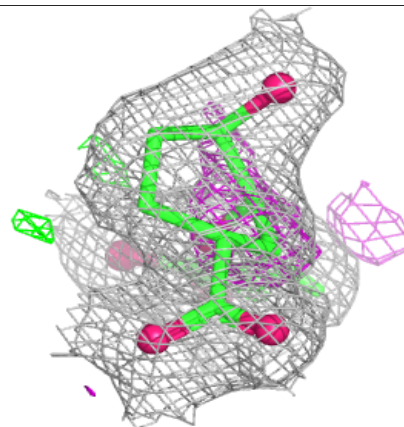
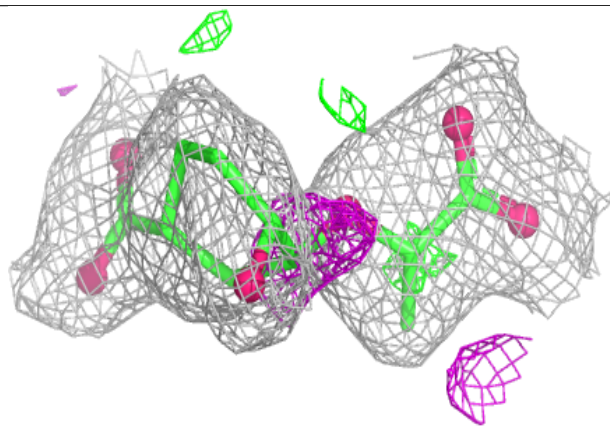
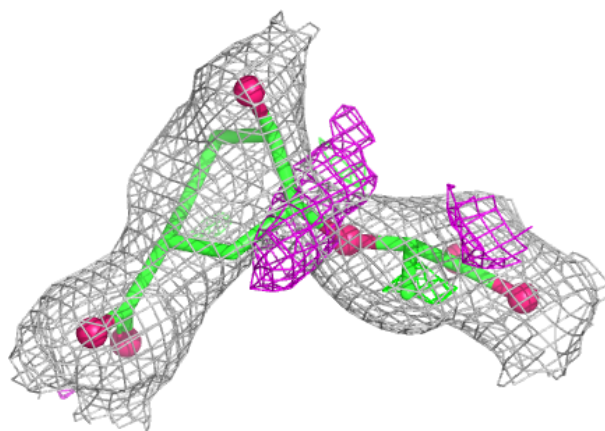
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FMT	A	603	3/3	0.89	0.14	39,39,43,44	0
3	MG	B	602	1/1	0.91	0.13	42,42,42,42	0
4	FMT	B	603	3/3	0.91	0.16	37,37,42,47	0
2	ISJ	B	601[B]	16/16	0.92	0.11	22,30,34,34	0
2	ISJ	A	601[B]	16/16	0.92	0.11	19,30,33,35	0
3	MG	A	602	1/1	0.98	0.08	42,42,42,42	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.

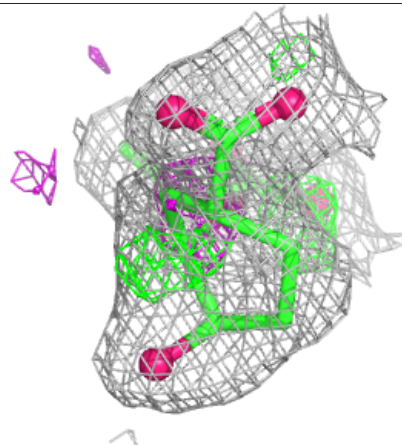
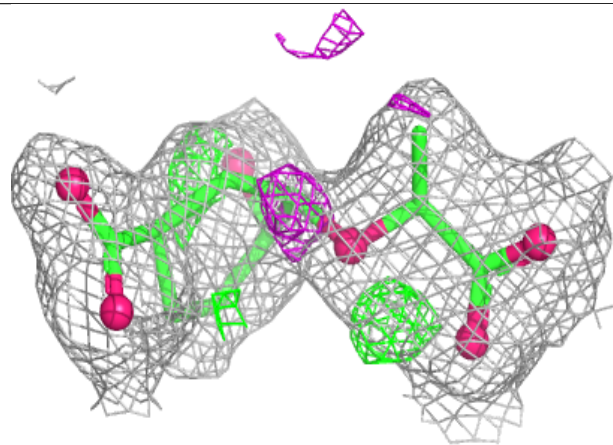
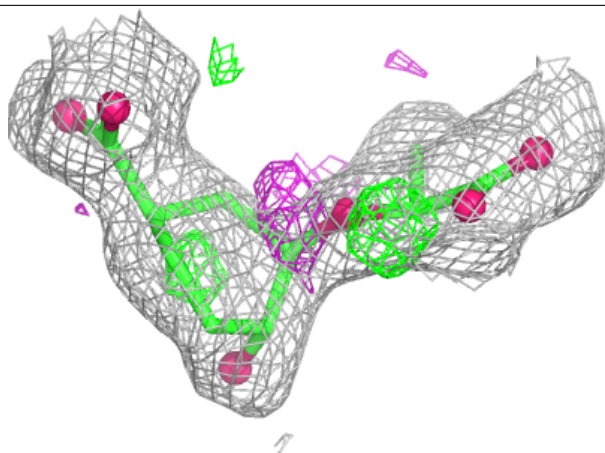
Electron density around ISJ B 601 (B):

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



Electron density around ISJ A 601 (B):

$2mF_o-DF_c$ (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.