



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 15, 2024 – 03:51 pm BST

PDB ID : 9GGJ
Title : Crystal structure of argininosuccinate lyase from Arabidopsis thaliana (AtASL) in complex with biological substrate and products - argininosuccinate, arginine and fumarate
Authors : Nielipinski, M.; Pietrzyk-Brzezinska, A.J.; Krzeszewska, D.; Sekula, B.
Deposited on : 2024-08-13
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
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

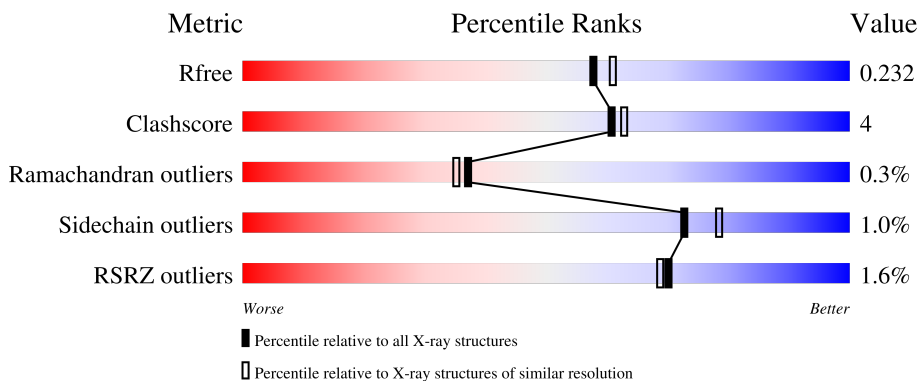
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.



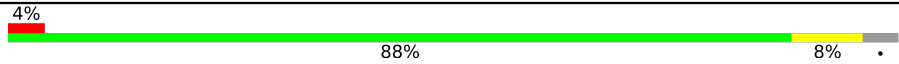
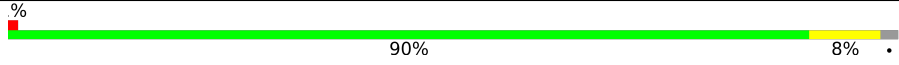
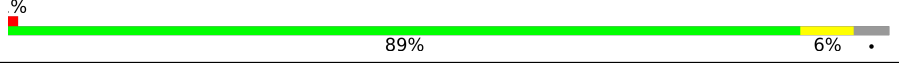

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

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 $\leq 5\%$. 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	465	 87% 10% .
1	B	465	 91% 6% .
1	C	465	 2% 86% 10% .
1	D	465	 87% 9% .

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Mol	Chain	Length	Quality of chain
1	E	465	 4% 88% 8% .
1	F	465	 % 90% 8% .
1	G	465	 % 89% 6% .
1	H	465	 2% 85% 11% ..

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 31810 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 Argininosuccinate lyase, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	451	Total 3572	C 2252	N 611	O 688	S 21	0	8	0
1	B	452	Total 3599	C 2271	N 616	O 691	S 21	0	10	0
1	C	448	Total 3555	C 2244	N 609	O 681	S 21	0	10	0
1	D	448	Total 3540	C 2235	N 607	O 678	S 20	0	6	0
1	E	448	Total 3552	C 2241	N 606	O 684	S 21	0	7	0
1	F	454	Total 3593	C 2265	N 618	O 690	S 20	0	6	0
1	G	448	Total 3548	C 2241	N 609	O 678	S 20	0	8	0
1	H	451	Total 3582	C 2264	N 610	O 687	S 21	0	8	0

There are 24 discrepancies between the modelled and reference sequences:

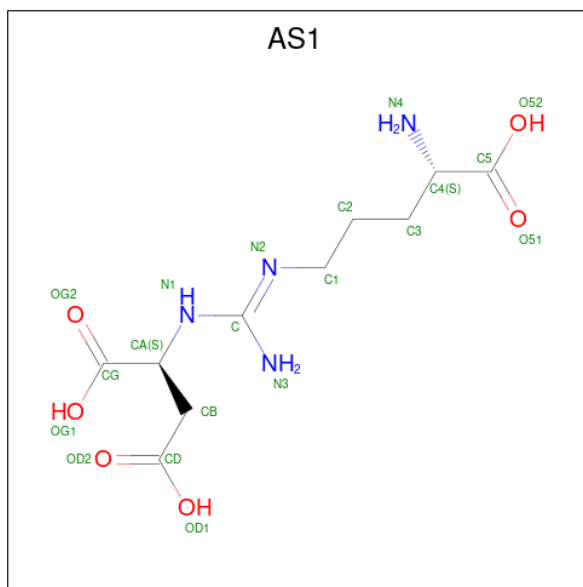
Chain	Residue	Modelled	Actual	Comment	Reference
A	53	SER	-	expression tag	UNP Q9LEU8
A	54	ASN	-	expression tag	UNP Q9LEU8
A	55	ALA	-	expression tag	UNP Q9LEU8
B	53	SER	-	expression tag	UNP Q9LEU8
B	54	ASN	-	expression tag	UNP Q9LEU8
B	55	ALA	-	expression tag	UNP Q9LEU8
C	53	SER	-	expression tag	UNP Q9LEU8
C	54	ASN	-	expression tag	UNP Q9LEU8
C	55	ALA	-	expression tag	UNP Q9LEU8
D	53	SER	-	expression tag	UNP Q9LEU8
D	54	ASN	-	expression tag	UNP Q9LEU8
D	55	ALA	-	expression tag	UNP Q9LEU8
E	53	SER	-	expression tag	UNP Q9LEU8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	54	ASN	-	expression tag	UNP Q9LEU8
E	55	ALA	-	expression tag	UNP Q9LEU8
F	53	SER	-	expression tag	UNP Q9LEU8
F	54	ASN	-	expression tag	UNP Q9LEU8
F	55	ALA	-	expression tag	UNP Q9LEU8
G	53	SER	-	expression tag	UNP Q9LEU8
G	54	ASN	-	expression tag	UNP Q9LEU8
G	55	ALA	-	expression tag	UNP Q9LEU8
H	53	SER	-	expression tag	UNP Q9LEU8
H	54	ASN	-	expression tag	UNP Q9LEU8
H	55	ALA	-	expression tag	UNP Q9LEU8

- Molecule 2 is ARGININOSUCCINATE (three-letter code: AS1) (formula: $C_{10}H_{18}N_4O_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			20	10	4	6		
2	D	1	Total	C	N	O	0	0
			20	10	4	6		
2	E	1	Total	C	N	O	0	0
			20	10	4	6		
2	G	1	Total	C	N	O	0	0
			20	10	4	6		
2	H	1	Total	C	N	O	0	0
			20	10	4	6		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

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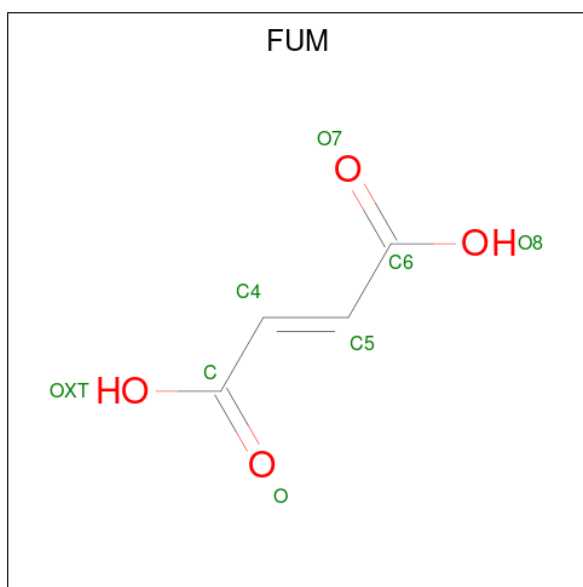
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
3	D	1	5	4	1	0	0
3	E	1	5	4	1	0	0
3	E	1	5	4	1	0	0
3	E	1	5	4	1	0	0
3	E	1	5	4	1	0	0
3	F	1	5	4	1	0	0
3	F	1	5	4	1	0	0
3	F	1	5	4	1	0	0
3	F	1	5	4	1	0	0
3	F	1	5	4	1	0	0
3	G	1	5	4	1	0	0
3	H	1	5	4	1	0	0
3	H	1	5	4	1	0	0
3	H	1	5	4	1	0	0
3	H	1	5	4	1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

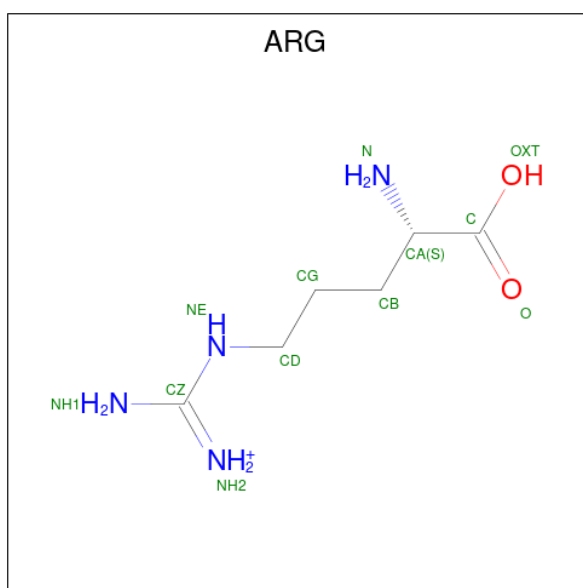
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
4	A	2	2	2	0	0
4	C	3	3	3	0	0
4	E	1	1	1	0	0
4	G	1	1	1	0	0

- Molecule 5 is FUMARIC ACID (three-letter code: FUM) (formula: C₄H₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	4	4		
5	E	1	Total	C	O	0	0
			8	4	4		

- Molecule 6 is ARGinine (three-letter code: ARG) (formula: $C_6H_{15}N_4O_2$).



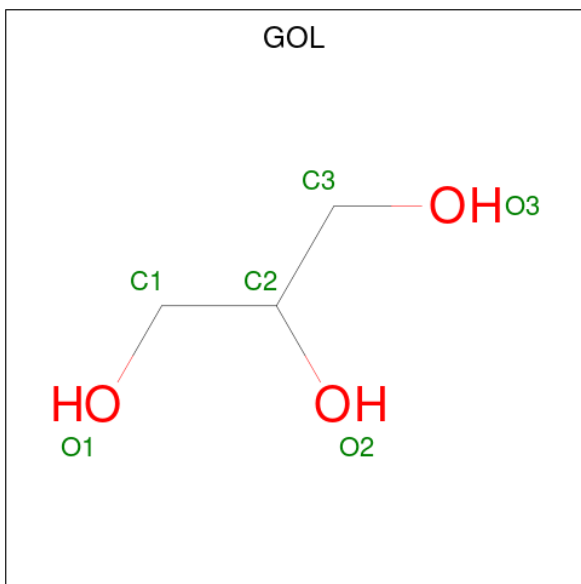
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			12	6	4	2		
6	C	1	Total	C	N	O	0	0
			12	6	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	F	1	12	6	4	2	0	0

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
7	D	1	6	3	3	0	0
7	G	1	6	3	3	0	0
7	G	1	6	3	3	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	353	353	353	0	0
8	B	344	344	344	0	0
8	C	376	376	376	0	0
8	D	400	400	400	0	0
8	E	331	331	331	0	0

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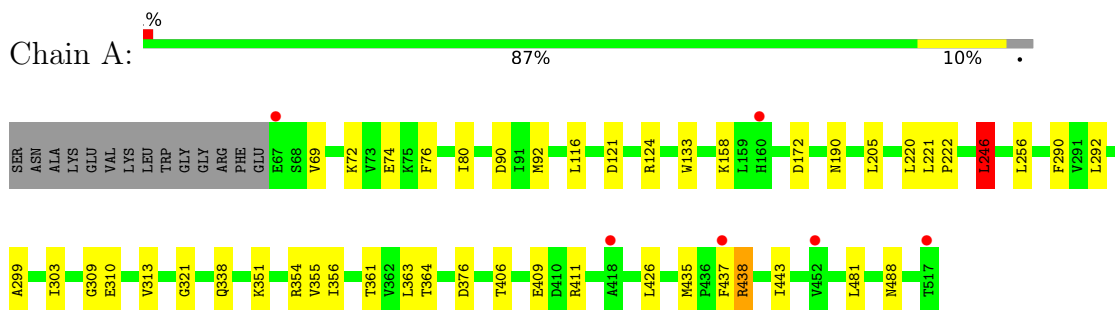
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	F	360	Total 360	O 360	0	0
8	G	407	Total 407	O 407	0	0
8	H	386	Total 386	O 386	0	0

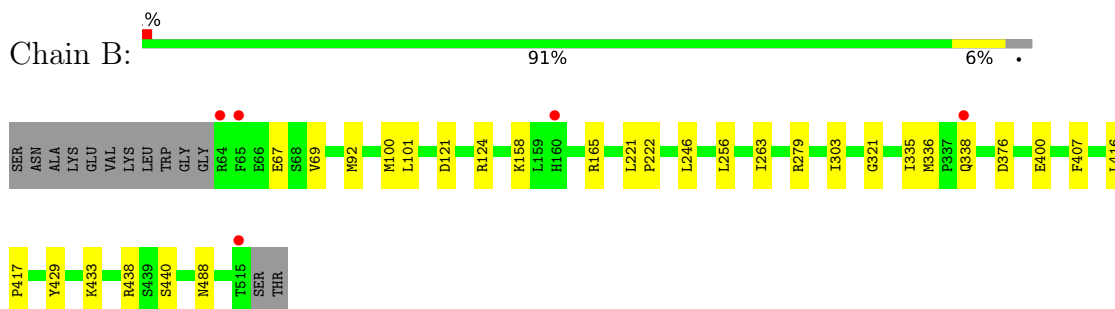
3 Residue-property plots

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.

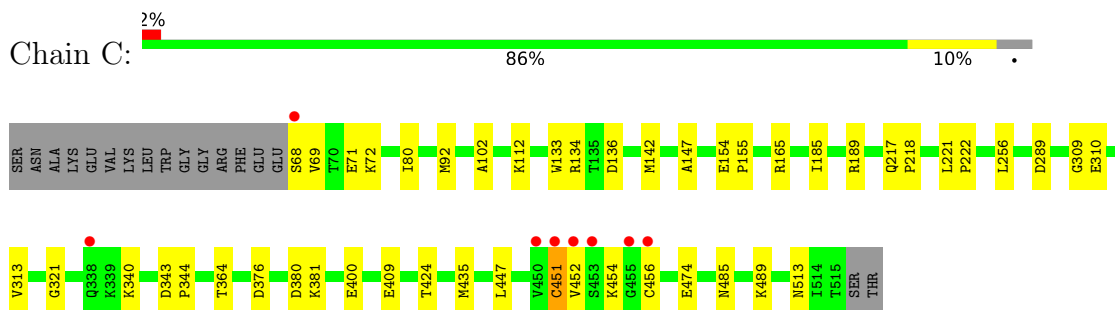
- Molecule 1: Argininosuccinate lyase, chloroplactic



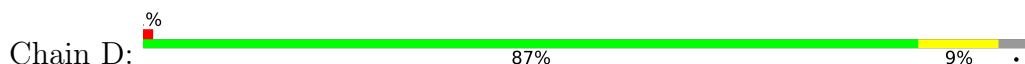
- Molecule 1: Argininosuccinate lyase, chloroplactic

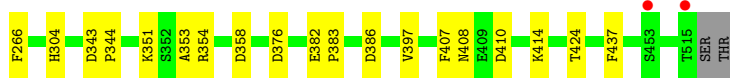


- Molecule 1: Argininosuccinate lyase, chloroplactic

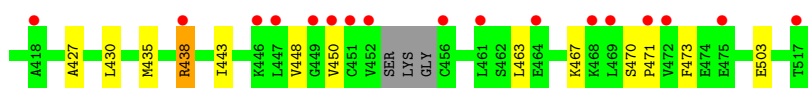
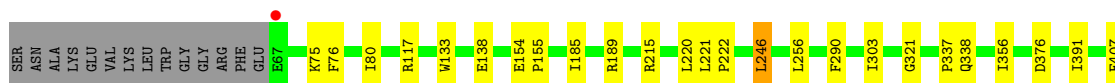
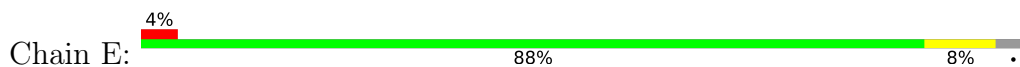


- Molecule 1: Argininosuccinate lyase, chloroplactic

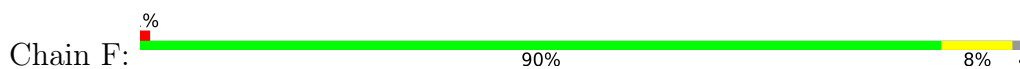




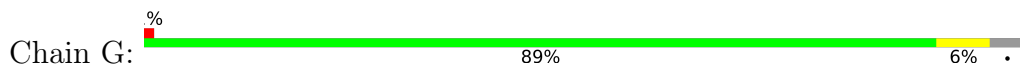
- Molecule 1: Argininosuccinate lyase, chloroplastic



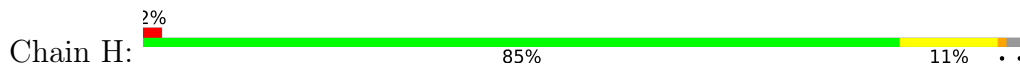
- Molecule 1: Argininosuccinate lyase, chloroplastic

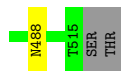


- Molecule 1: Argininosuccinate lyase, chloroplastic



- Molecule 1: Argininosuccinate lyase, chloroplastic





4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.59Å 230.85Å 112.53Å 90.00° 89.48° 90.00°	Depositor
Resolution (Å)	49.57 – 2.00 49.57 – 2.00	Depositor EDS
% Data completeness (in resolution range)	71.6 (49.57-2.00) 71.6 (49.57-2.00)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.00Å)	Xtrriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.175 , 0.227 0.186 , 0.232	Depositor DCC
R_{free} test set	258808 reflections (0.76%)	wwPDB-VP
Wilson B-factor (Å ²)	21.0	Xtrriage
Anisotropy	0.356	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	31810	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: GOL, CL, FUM, AS1, SO4

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	0.49	0/3659	0.83	1/4946 (0.0%)
1	B	0.49	0/3693	0.85	0/4992
1	C	0.52	0/3648	0.89	0/4932
1	D	0.51	0/3621	0.86	0/4896
1	E	0.48	0/3632	0.85	0/4910
1	F	0.51	0/3675	0.84	0/4965
1	G	0.54	0/3635	0.87	0/4913
1	H	0.50	0/3670	0.88	1/4960 (0.0%)
All	All	0.50	0/29233	0.86	2/39514 (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
1	H	0	1
All	All	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	LEU	N-CA-CB	-7.10	96.19	110.40
1	H	137	ARG	CB-CA-C	5.08	120.56	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	438	ARG	Sidechain
1	H	411	ARG	Sidechain

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	3572	0	3576	38	0
1	B	3599	0	3605	27	0
1	C	3555	0	3573	39	0
1	D	3540	0	3550	31	0
1	E	3552	0	3550	29	0
1	F	3593	0	3589	29	0
1	G	3548	0	3568	25	0
1	H	3582	0	3590	38	0
2	A	20	0	14	1	0
2	D	20	0	14	0	0
2	E	20	0	14	1	0
2	G	20	0	14	0	0
2	H	20	0	14	1	0
3	A	25	0	0	1	0
3	B	20	0	0	1	0
3	C	15	0	0	0	0
3	D	10	0	0	1	0
3	E	20	0	0	0	0
3	F	20	0	0	2	0
3	G	5	0	0	0	0
3	H	20	0	0	0	0
4	A	2	0	0	0	0
4	C	3	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
5	A	8	0	1	0	0
5	E	8	0	1	0	0
6	B	12	0	12	1	0
6	C	12	0	12	1	0
6	F	12	0	12	0	0
7	D	6	0	8	0	0
7	G	12	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	353	0	0	11	0
8	B	344	0	0	9	0
8	C	376	0	0	18	0
8	D	400	0	0	10	0
8	E	331	0	0	7	0
8	F	360	0	0	8	0
8	G	407	0	0	5	0
8	H	386	0	0	9	0
All	All	31810	0	28733	235	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 4.

The worst 5 of 235 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:364[B]:THR:HG22	1:C:364[B]:THR:HG22	1.54	0.88
1:H:318:GLU:HB2	8:H:855:HOH:O	1.77	0.85
1:A:190:ASN:HB2	8:A:900:HOH:O	1.79	0.81
1:A:338:GLN:HB3	8:A:994:HOH:O	1.81	0.79
1:B:335:ILE:HG13	1:B:336:MET:HG3	1.62	0.79

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	457/465 (98%)	445 (97%)	11 (2%)	1 (0%)	44 42
1	B	460/465 (99%)	452 (98%)	7 (2%)	1 (0%)	44 42
1	C	456/465 (98%)	443 (97%)	12 (3%)	1 (0%)	44 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	452/465 (97%)	444 (98%)	7 (2%)	1 (0%)	44	42
1	E	451/465 (97%)	439 (97%)	10 (2%)	2 (0%)	30	27
1	F	458/465 (98%)	447 (98%)	10 (2%)	1 (0%)	44	42
1	G	454/465 (98%)	445 (98%)	8 (2%)	1 (0%)	44	42
1	H	457/465 (98%)	442 (97%)	12 (3%)	3 (1%)	19	14
All	All	3645/3720 (98%)	3557 (98%)	77 (2%)	11 (0%)	37	35

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	256	LEU
1	B	256	LEU
1	E	256	LEU
1	C	256	LEU
1	D	256	LEU

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	399/402 (99%)	393 (98%)	6 (2%)	60	66
1	B	402/402 (100%)	400 (100%)	2 (0%)	86	90
1	C	398/402 (99%)	396 (100%)	2 (0%)	86	90
1	D	394/402 (98%)	390 (99%)	4 (1%)	73	78
1	E	396/402 (98%)	393 (99%)	3 (1%)	79	84
1	F	398/402 (99%)	394 (99%)	4 (1%)	73	78
1	G	396/402 (98%)	392 (99%)	4 (1%)	73	78
1	H	399/402 (99%)	392 (98%)	7 (2%)	54	59
All	All	3182/3216 (99%)	3150 (99%)	32 (1%)	73	78

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	376	ASP
1	H	432	LYS
1	D	376	ASP
1	D	246	LEU
1	H	451	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	300	ASN
1	A	338	GLN
1	F	214	GLN
1	F	300	ASN
1	G	300	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 47 ligands modelled in this entry, 7 are monoatomic - leaving 40 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
7	GOL	G	602	-	5,5,5	0.08	0	5,5,5	0.36	0
3	SO4	A	602	-	4,4,4	0.44	0	6,6,6	0.29	0
3	SO4	D	604	-	4,4,4	0.28	0	6,6,6	0.18	0
3	SO4	E	605	-	4,4,4	0.28	0	6,6,6	0.35	0
6	ARG	C	707	-	10,11,11	0.61	0	11,13,13	0.98	0
2	AS1	G	601	-	18,19,19	1.13	2 (11%)	21,24,24	1.12	2 (9%)
3	SO4	F	703	-	4,4,4	0.32	0	6,6,6	0.13	0
6	ARG	B	605	-	10,11,11	0.65	0	11,13,13	0.79	0
2	AS1	D	601	-	18,19,19	1.12	2 (11%)	21,24,24	1.07	1 (4%)
3	SO4	E	604	-	4,4,4	0.37	0	6,6,6	0.30	0
3	SO4	H	604	-	4,4,4	0.34	0	6,6,6	0.10	0
3	SO4	A	604	-	4,4,4	0.30	0	6,6,6	0.39	0
5	FUM	A	609	-	7,7,7	1.91	3 (42%)	8,8,8	0.33	0
3	SO4	E	603	-	4,4,4	0.36	0	6,6,6	0.05	0
2	AS1	H	601	-	18,19,19	1.05	1 (5%)	21,24,24	1.16	1 (4%)
3	SO4	D	603	-	4,4,4	0.36	0	6,6,6	0.27	0
3	SO4	F	701	-	4,4,4	0.36	0	6,6,6	0.17	0
3	SO4	G	604	-	4,4,4	0.31	0	6,6,6	0.17	0
7	GOL	D	602	-	5,5,5	0.19	0	5,5,5	0.43	0
6	ARG	F	705	-	10,11,11	0.78	0	11,13,13	1.00	1 (9%)
2	AS1	A	601	-	18,19,19	1.08	0	21,24,24	1.11	1 (4%)
3	SO4	A	603	-	4,4,4	0.29	0	6,6,6	0.14	0
3	SO4	C	702	-	4,4,4	0.34	0	6,6,6	0.18	0
3	SO4	E	606	-	4,4,4	0.28	0	6,6,6	0.15	0
3	SO4	H	603	-	4,4,4	0.38	0	6,6,6	0.24	0
7	GOL	G	603	-	5,5,5	0.21	0	5,5,5	0.43	0
3	SO4	F	704	-	4,4,4	0.34	0	6,6,6	0.22	0
3	SO4	H	605	-	4,4,4	0.38	0	6,6,6	0.11	0
3	SO4	B	601	-	4,4,4	0.40	0	6,6,6	0.19	0
3	SO4	A	605	-	4,4,4	0.32	0	6,6,6	0.13	0
3	SO4	C	703	-	4,4,4	0.38	0	6,6,6	0.19	0
3	SO4	B	602	-	4,4,4	0.37	0	6,6,6	0.22	0
3	SO4	B	604	-	4,4,4	0.34	0	6,6,6	0.10	0
3	SO4	C	701	-	4,4,4	0.38	0	6,6,6	0.28	0
5	FUM	E	602	-	7,7,7	1.78	2 (28%)	8,8,8	0.39	0
2	AS1	E	601	-	18,19,19	0.99	2 (11%)	21,24,24	1.17	2 (9%)
3	SO4	B	603	-	4,4,4	0.29	0	6,6,6	0.29	0
3	SO4	F	702	-	4,4,4	0.30	0	6,6,6	0.35	0
3	SO4	H	602	-	4,4,4	0.44	0	6,6,6	0.20	0
3	SO4	A	606	-	4,4,4	0.35	0	6,6,6	0.14	0

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
7	GOL	G	602	-	-	3/4/4/4	-
5	FUM	E	602	-	-	0/5/5/5	-
2	AS1	H	601	-	-	2/21/23/23	-
2	AS1	E	601	-	-	3/21/23/23	-
7	GOL	D	602	-	-	0/4/4/4	-
6	ARG	C	707	-	-	1/11/11/11	-
2	AS1	A	601	-	-	4/21/23/23	-
2	AS1	G	601	-	-	3/21/23/23	-
6	ARG	F	705	-	-	3/11/11/11	-
6	ARG	B	605	-	-	3/11/11/11	-
2	AS1	D	601	-	-	3/21/23/23	-
7	GOL	G	603	-	-	2/4/4/4	-
5	FUM	A	609	-	-	2/5/5/5	-

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	609	FUM	O8-C6	-3.29	1.21	1.30
5	E	602	FUM	OXT-C	-3.10	1.22	1.30
2	D	601	AS1	O51-C5	2.72	1.30	1.22
5	E	602	FUM	O8-C6	-2.67	1.23	1.30
2	E	601	AS1	OD1-CD	-2.34	1.22	1.30

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	601	AS1	C2-C1-N2	2.64	115.46	110.66
2	D	601	AS1	C2-C1-N2	2.44	115.10	110.66
2	G	601	AS1	N1-C-N2	-2.38	119.10	124.90
2	E	601	AS1	N1-C-N2	-2.31	119.27	124.90
6	F	705	ARG	CB-CA-C	2.27	115.70	110.30

There are no chirality outliers.

5 of 29 torsion outliers are listed below:

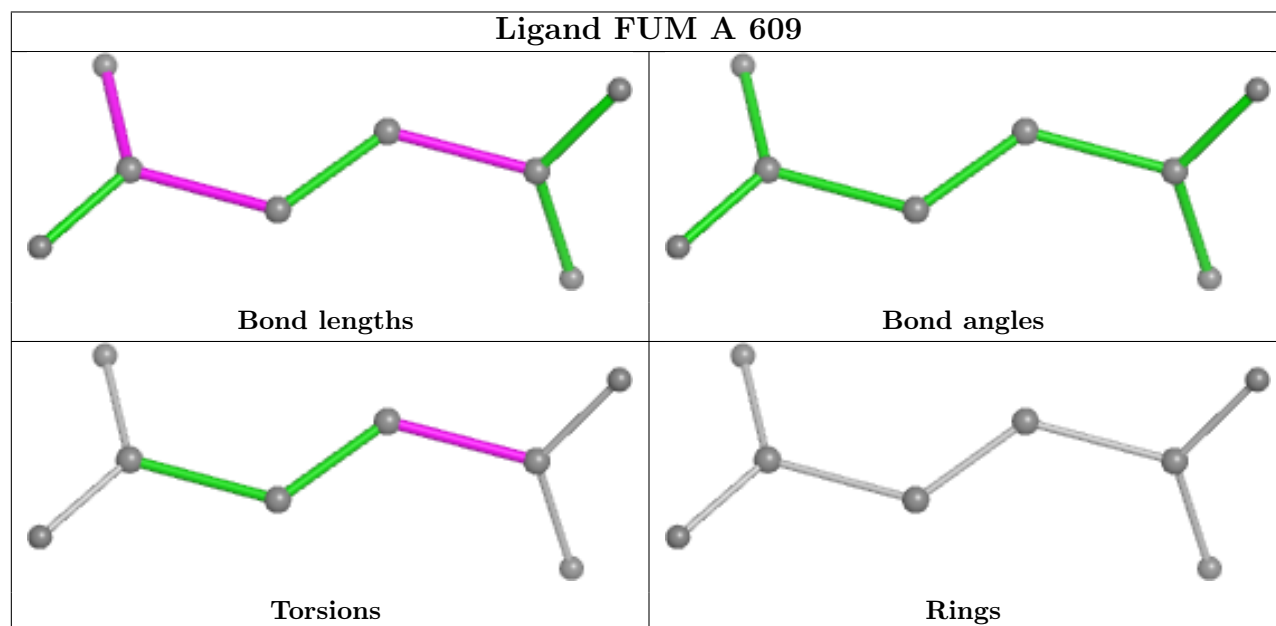
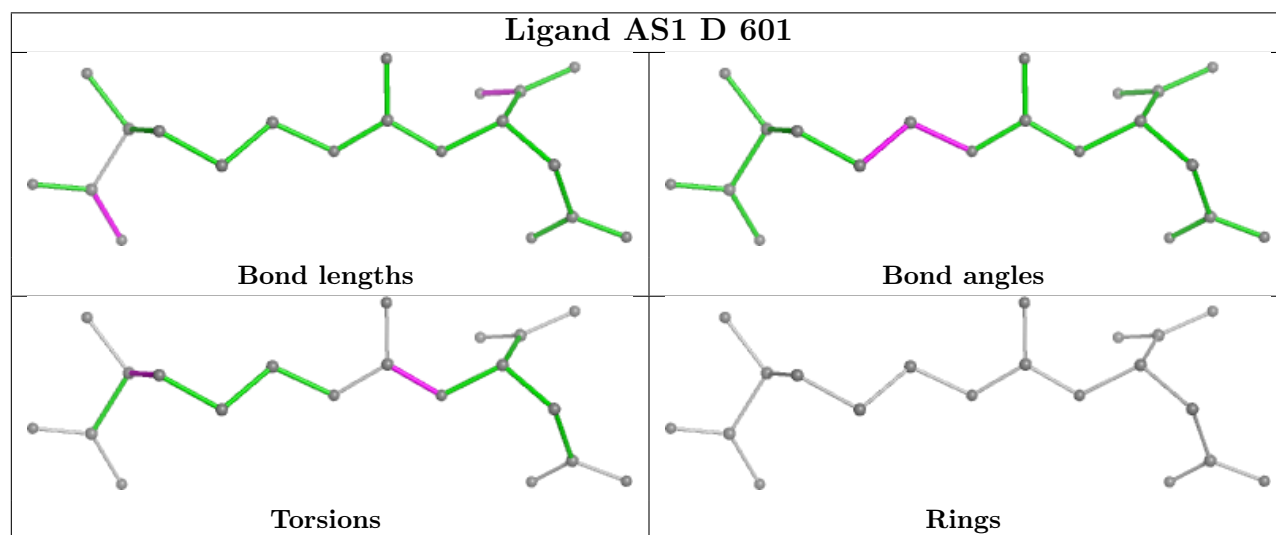
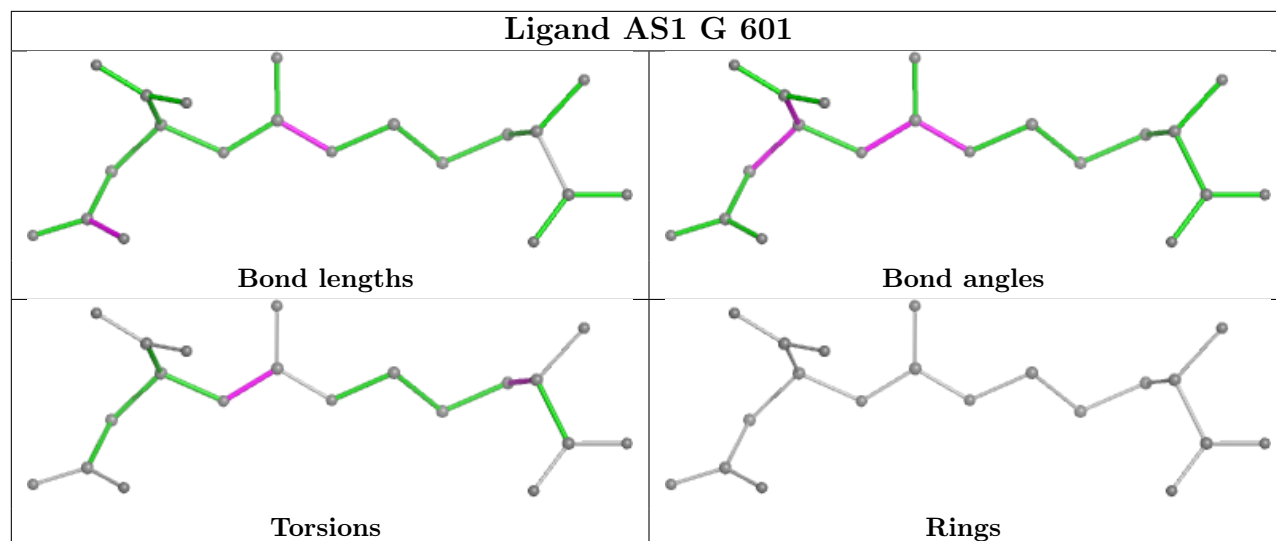
Mol	Chain	Res	Type	Atoms
2	A	601	AS1	C2-C3-C4-N4
2	D	601	AS1	N2-C-N1-CA
2	G	601	AS1	N2-C-N1-CA
2	H	601	AS1	N2-C-N1-CA
6	B	605	ARG	N-CA-CB-CG

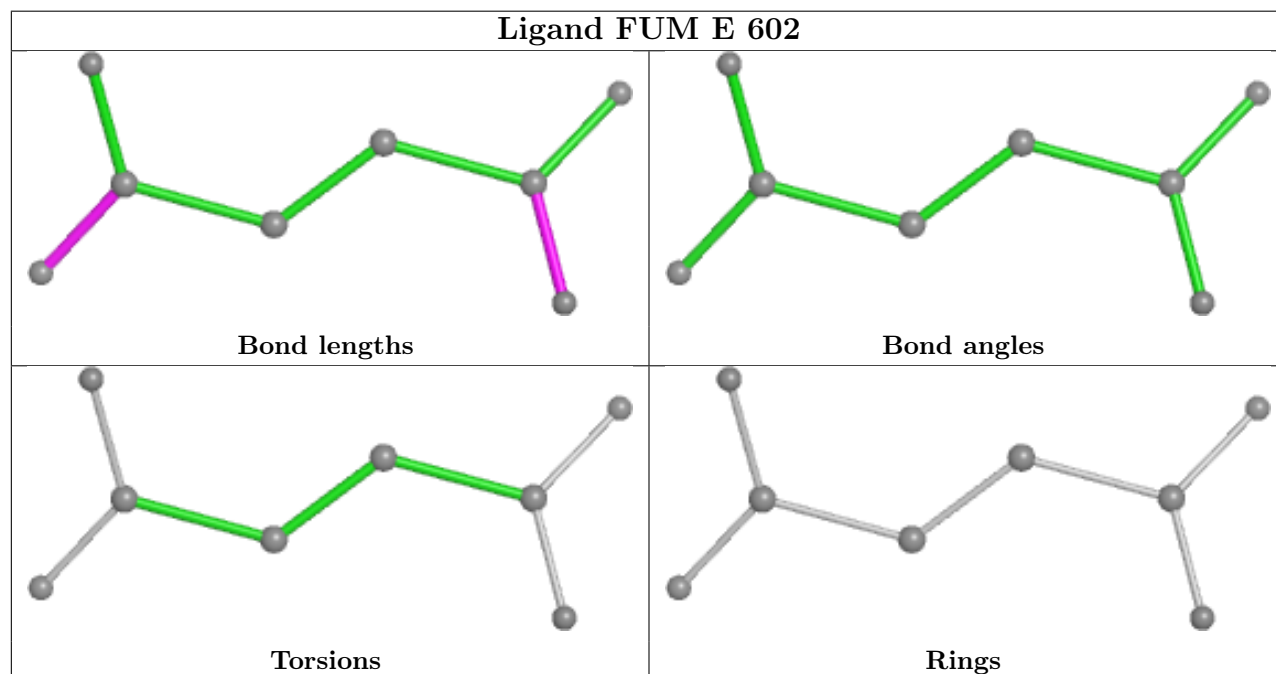
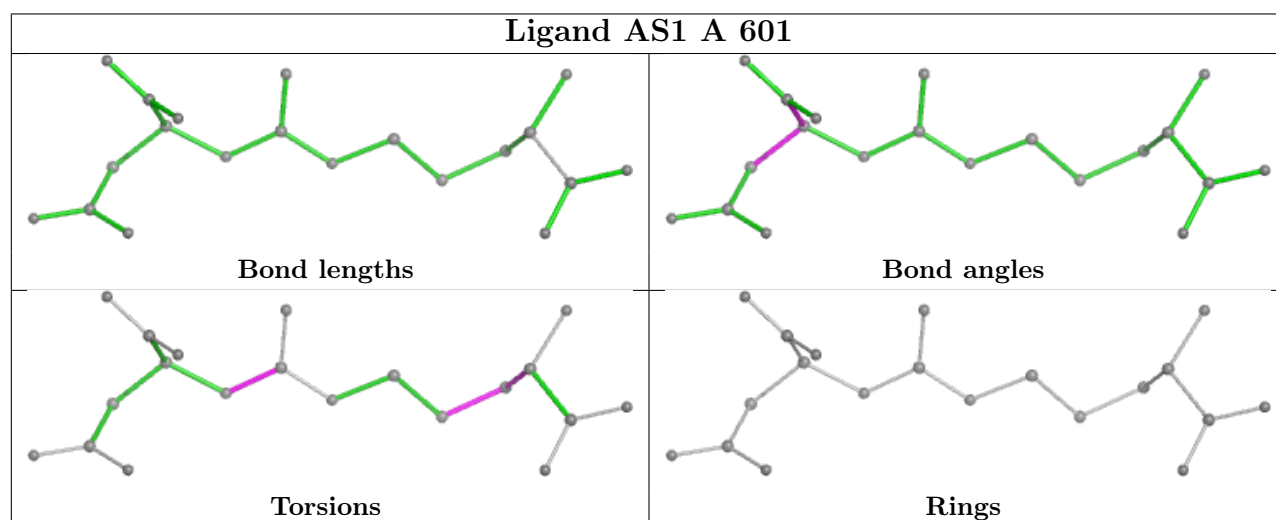
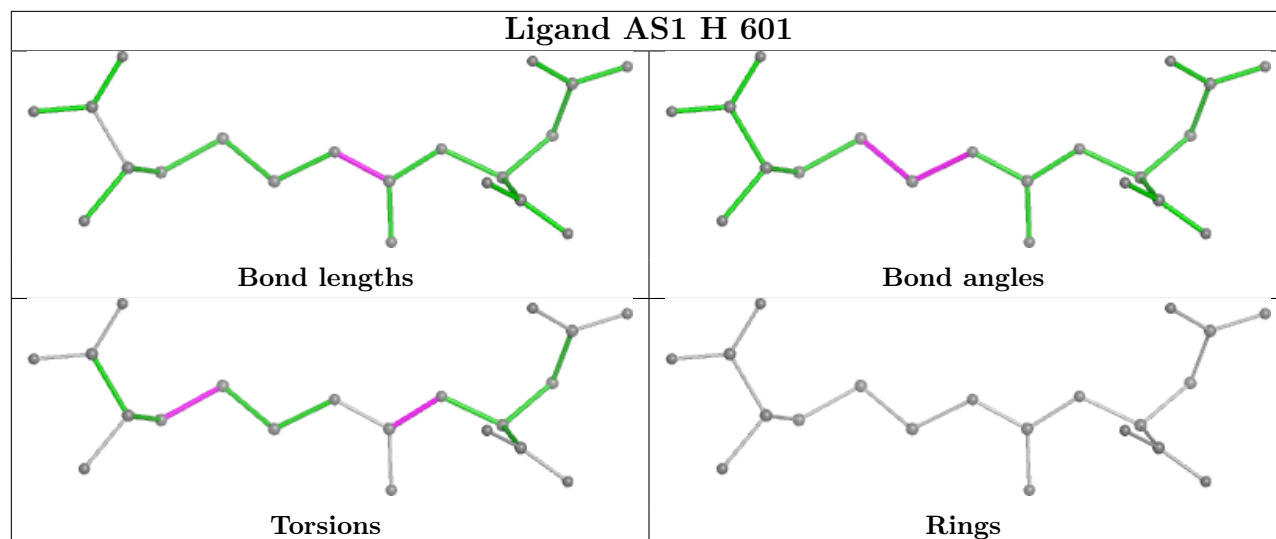
There are no ring outliers.

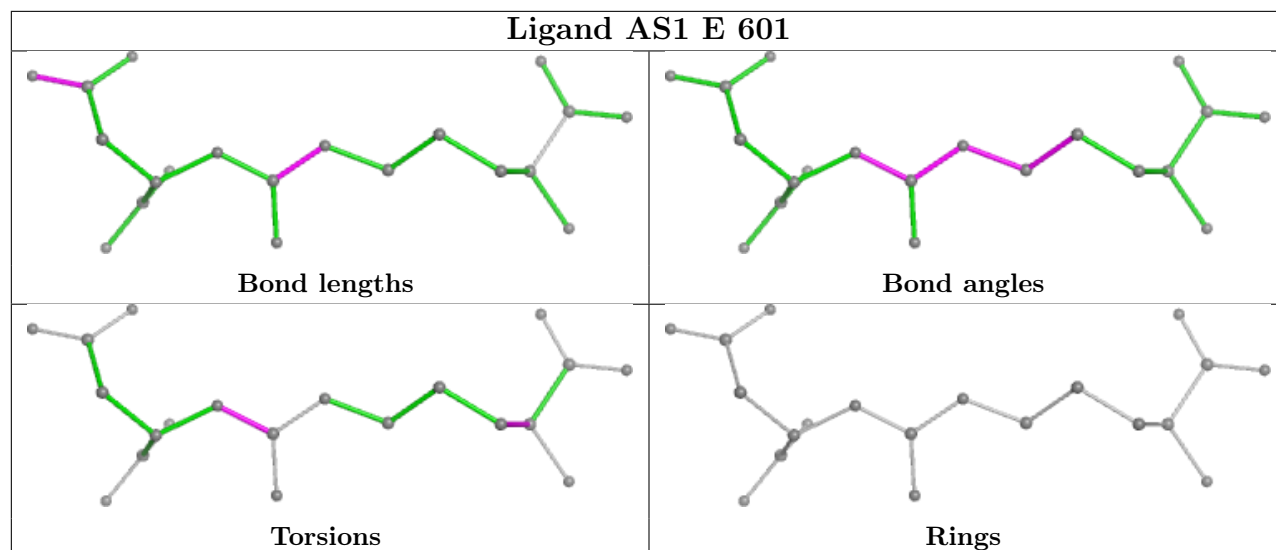
11 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	G	602	GOL	1	0
3	A	602	SO4	1	0
3	D	604	SO4	1	0
6	C	707	ARG	1	0
6	B	605	ARG	1	0
2	H	601	AS1	1	0
2	A	601	AS1	1	0
3	F	704	SO4	1	0
3	B	601	SO4	1	0
2	E	601	AS1	1	0
3	F	702	SO4	1	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	451/465 (96%)	-0.27	6 (1%) 74 73	12, 24, 50, 81	8 (1%)
1	B	452/465 (97%)	-0.33	5 (1%) 77 76	11, 24, 45, 83	10 (2%)
1	C	448/465 (96%)	-0.23	8 (1%) 67 66	11, 23, 52, 97	10 (2%)
1	D	448/465 (96%)	-0.32	3 (0%) 84 83	11, 23, 44, 61	6 (1%)
1	E	448/465 (96%)	-0.08	18 (4%) 43 41	12, 24, 67, 93	7 (1%)
1	F	454/465 (97%)	-0.27	4 (0%) 81 80	12, 25, 47, 91	6 (1%)
1	G	448/465 (96%)	-0.33	3 (0%) 84 83	10, 22, 46, 81	8 (1%)
1	H	451/465 (96%)	-0.28	11 (2%) 59 58	11, 23, 49, 95	8 (1%)
All	All	3600/3720 (96%)	-0.27	58 (1%) 70 69	10, 24, 51, 97	63 (1%)

The worst 5 of 58 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	452	VAL	5.6
1	E	418	ALA	4.4
1	H	455	GLY	4.2
1	G	337	PRO	4.1
1	A	437	PHE	4.0

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

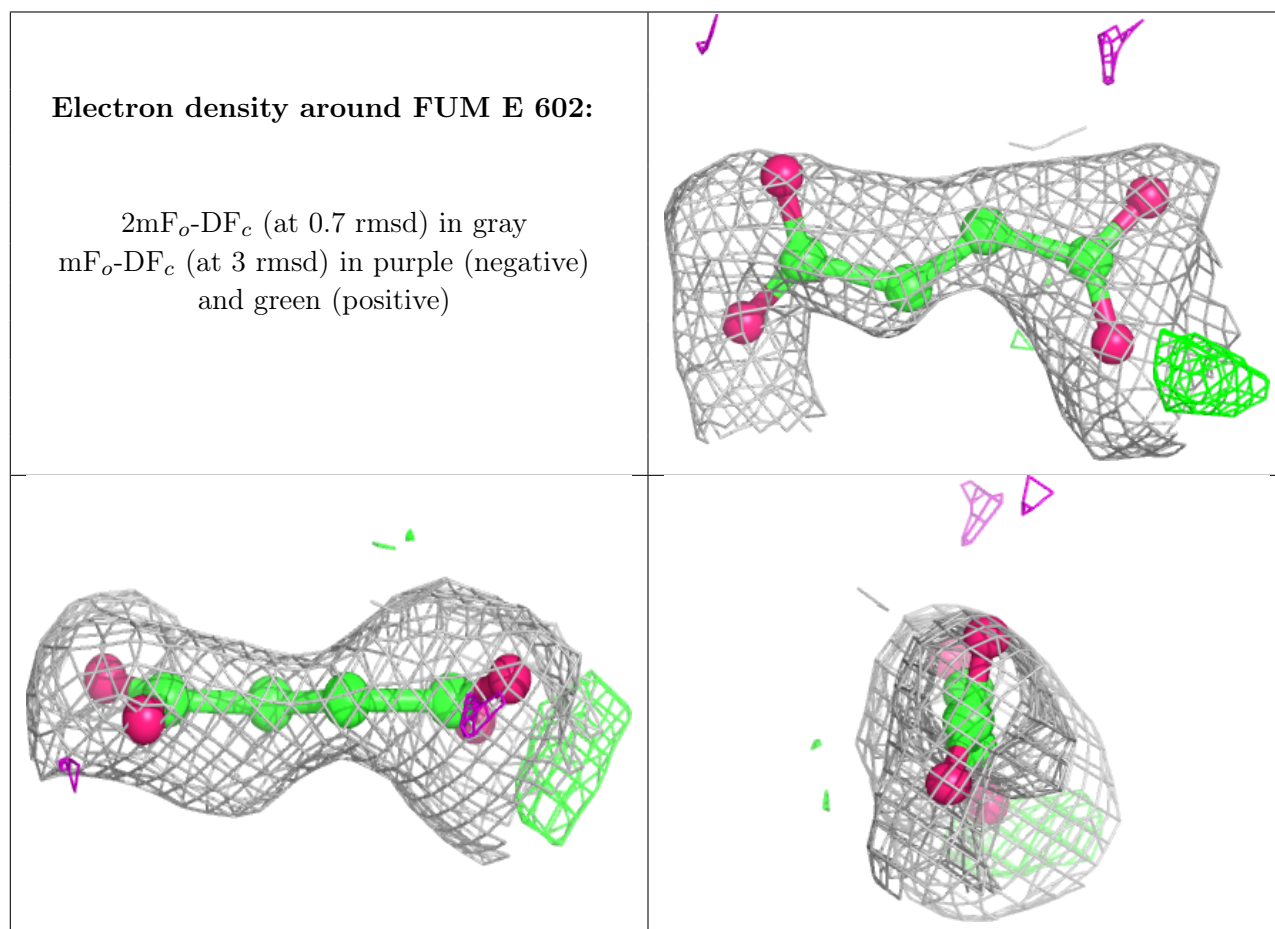
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	E	606	5/5	0.76	0.16	76,77,98,98	0
7	GOL	G	602	6/6	0.77	0.19	42,47,50,52	0
3	SO4	C	703	5/5	0.80	0.15	69,82,87,89	0
3	SO4	B	604	5/5	0.81	0.12	75,82,86,93	0
3	SO4	F	704	5/5	0.82	0.14	73,78,84,87	0
3	SO4	A	605	5/5	0.82	0.12	73,74,80,81	0
3	SO4	H	603	5/5	0.87	0.13	44,46,61,70	0
7	GOL	D	602	6/6	0.87	0.12	49,52,53,54	0
3	SO4	B	602	5/5	0.87	0.12	64,68,82,94	0
5	FUM	E	602	8/8	0.88	0.15	42,55,71,71	0
3	SO4	E	603	5/5	0.88	0.12	81,84,93,96	0
3	SO4	F	703	5/5	0.88	0.08	71,79,81,81	0
7	GOL	G	603	6/6	0.88	0.15	47,53,56,57	0
3	SO4	H	604	5/5	0.89	0.12	63,88,91,96	0
3	SO4	H	605	5/5	0.89	0.14	69,97,105,111	0
3	SO4	B	603	5/5	0.89	0.12	53,57,65,85	0
3	SO4	C	702	5/5	0.90	0.12	58,69,76,83	0
2	AS1	H	601	20/20	0.90	0.12	33,50,60,63	0
6	ARG	C	707	12/12	0.90	0.10	23,30,39,45	0
3	SO4	A	606	5/5	0.91	0.10	75,75,78,82	0
3	SO4	A	602	5/5	0.91	0.10	49,53,56,58	0
6	ARG	F	705	12/12	0.91	0.11	31,38,49,51	0
3	SO4	F	702	5/5	0.91	0.12	49,58,62,63	0
2	AS1	E	601	20/20	0.91	0.10	31,43,50,51	0
5	FUM	A	609	8/8	0.91	0.13	40,58,62,65	0
3	SO4	E	604	5/5	0.92	0.13	47,56,66,81	0
3	SO4	E	605	5/5	0.92	0.13	39,40,52,59	0
3	SO4	A	604	5/5	0.92	0.15	42,43,51,61	0
3	SO4	B	601	5/5	0.92	0.10	55,56,65,71	0
3	SO4	D	604	5/5	0.93	0.09	53,54,63,77	0
2	AS1	D	601	20/20	0.93	0.10	22,41,54,55	0
2	AS1	A	601	20/20	0.93	0.10	30,43,56,58	0
3	SO4	A	603	5/5	0.93	0.10	41,47,57,63	0
3	SO4	H	602	5/5	0.93	0.10	54,56,57,73	0
3	SO4	D	603	5/5	0.94	0.11	44,57,63,78	0
2	AS1	G	601	20/20	0.94	0.09	21,33,45,46	0
6	ARG	B	605	12/12	0.95	0.07	25,33,36,37	0

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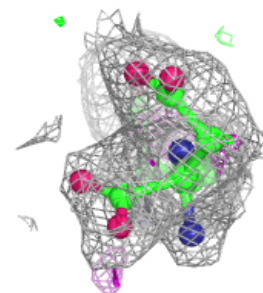
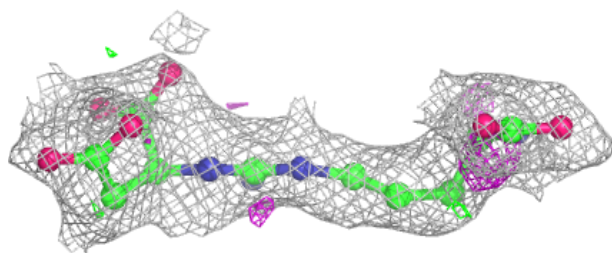
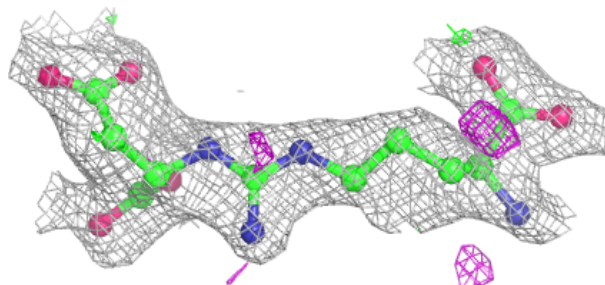
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	C	701	5/5	0.96	0.07	50,53,58,60	0
3	SO4	G	604	5/5	0.96	0.12	40,42,50,53	0
3	SO4	F	701	5/5	0.97	0.13	42,43,49,54	0
4	CL	E	607	1/1	0.98	0.05	32,32,32,32	0
4	CL	G	605	1/1	0.98	0.07	23,23,23,23	0
4	CL	A	608	1/1	0.98	0.08	31,31,31,31	0
4	CL	C	705	1/1	0.98	0.05	26,26,26,26	0
4	CL	C	706	1/1	0.98	0.07	22,22,22,22	0
4	CL	A	607	1/1	0.99	0.10	34,34,34,34	0
4	CL	C	704	1/1	0.99	0.05	22,22,22,22	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.

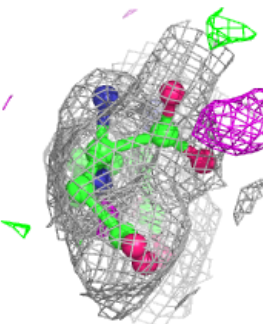
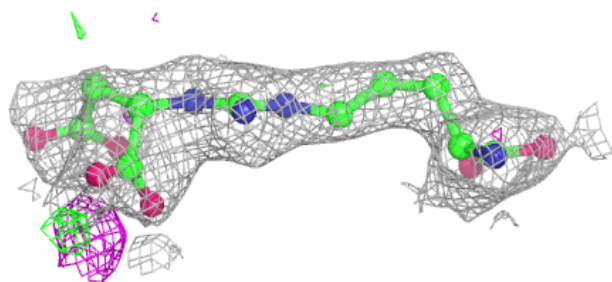
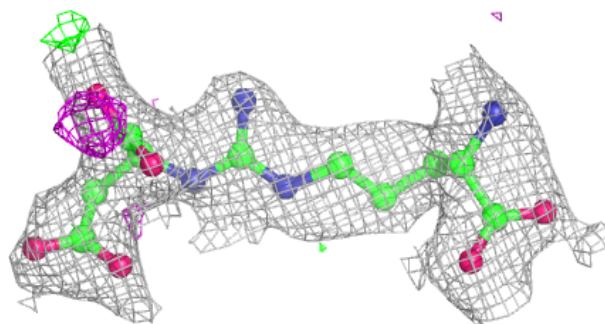


Electron density around AS1 H 601:

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

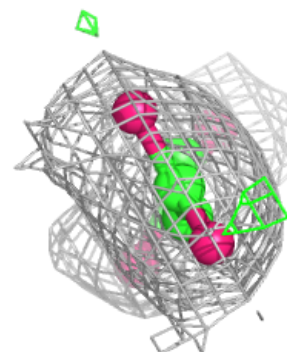
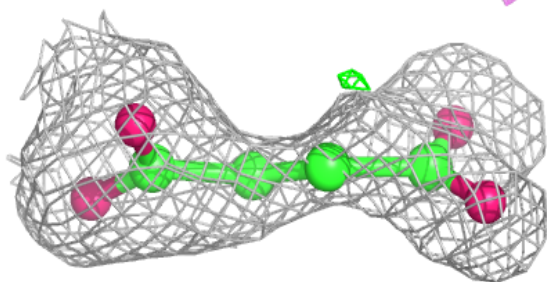
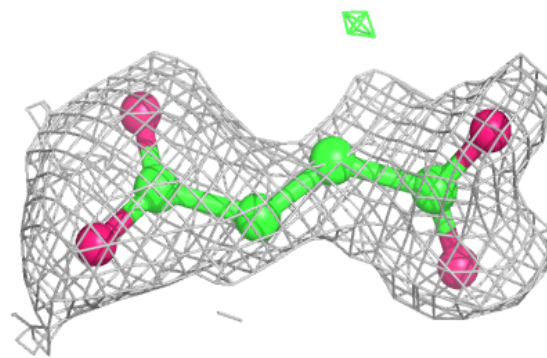
**Electron density around AS1 E 601:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

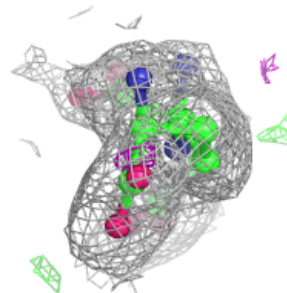
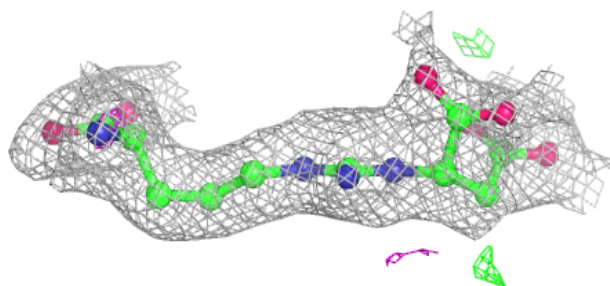
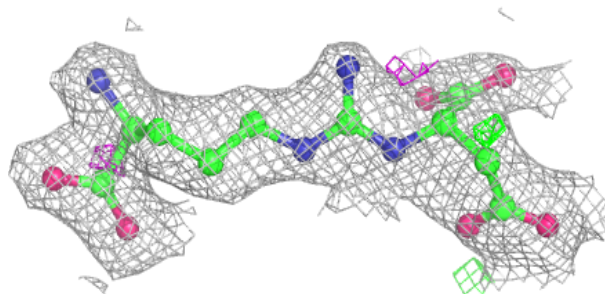


Electron density around FUM A 609:

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

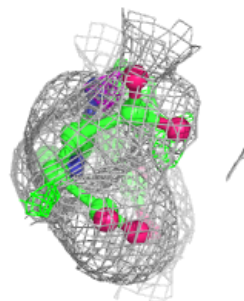
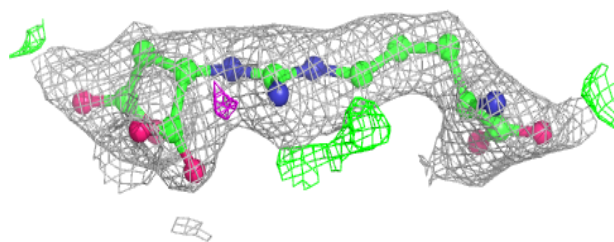
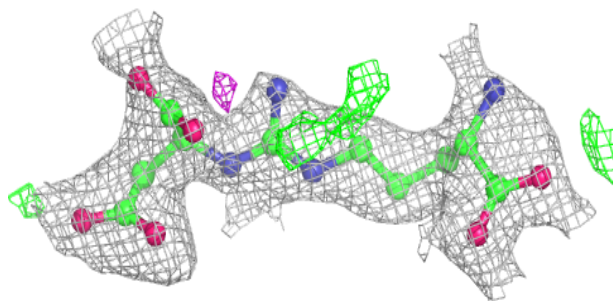
**Electron density around AS1 D 601:**

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

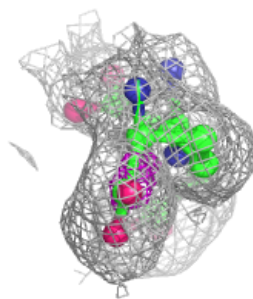
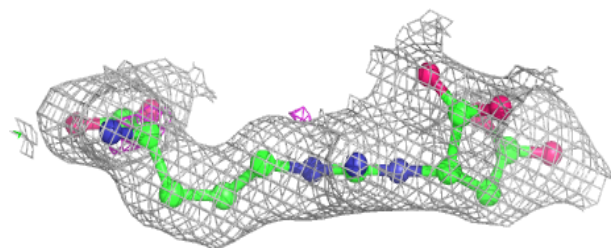
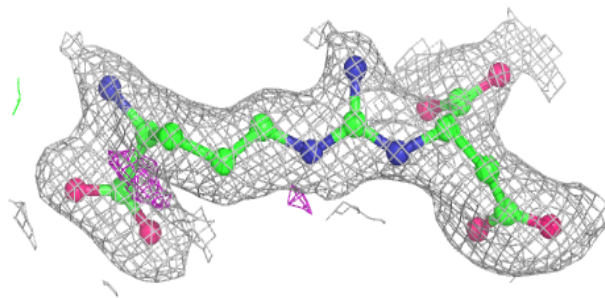


Electron density around AS1 A 601:

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

**Electron density around AS1 G 601:**

$2mF_o-DF_c$ (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.