



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 10:09 PM BST

PDB ID : 6O4I
Title : Structure of ALDH7A1 mutant E399D complexed with alpha-aminoadipate
Authors : Tanner, J.J.; Korasick, D.A.; Laciak, A.R.
Deposited on : 2019-02-28
Resolution : 1.75 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

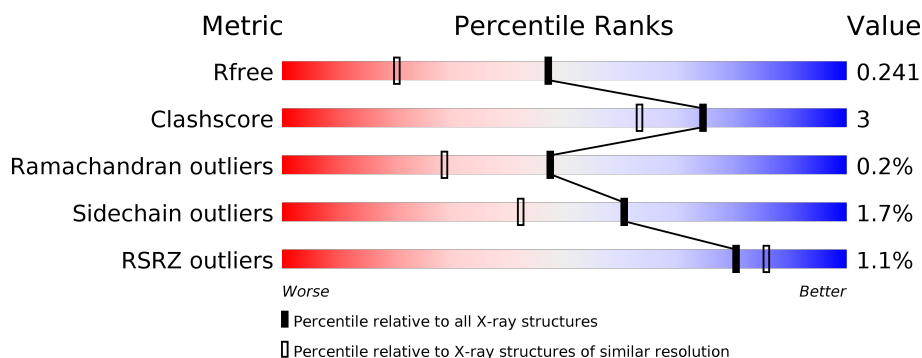
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13
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) : 2.13

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X-RAY DIFFRACTION

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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

Mol	Chain	Length	Quality of chain
1	A	513	<div> <div></div> <div>91%</div> <div>9%</div> </div>
1	B	513	<div> <div></div> <div>88%</div> <div>11%</div> </div>
1	C	513	<div> <div></div> <div>90%</div> <div>9%</div> </div>
1	D	513	<div> <div></div> <div>90%</div> <div>9%</div> </div>
1	E	513	<div> <div></div> <div>90%</div> <div>9%</div> </div>
1	F	513	<div> <div></div> <div>89%</div> <div>10%</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	513	<div><div>%</div><div><div></div></div><div>88%11%•</div></div>
1	H	513	<div><div>%</div><div><div></div></div><div>88%11%•</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 33043 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 Alpha-aminoadipic semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	509	Total	C	N	O	S	0	3	0
			3860	2449	668	725	18			
1	B	509	Total	C	N	O	S	0	2	0
			3866	2455	670	723	18			
1	C	509	Total	C	N	O	S	0	1	0
			3866	2454	671	723	18			
1	D	509	Total	C	N	O	S	0	1	0
			3864	2453	668	725	18			
1	E	509	Total	C	N	O	S	0	2	0
			3881	2464	674	725	18			
1	F	509	Total	C	N	O	S	0	1	0
			3878	2462	674	724	18			
1	G	509	Total	C	N	O	S	0	2	0
			3871	2460	672	721	18			
1	H	509	Total	C	N	O	S	0	1	0
			3870	2457	671	724	18			

There are 24 discrepancies between the modelled and reference sequences:

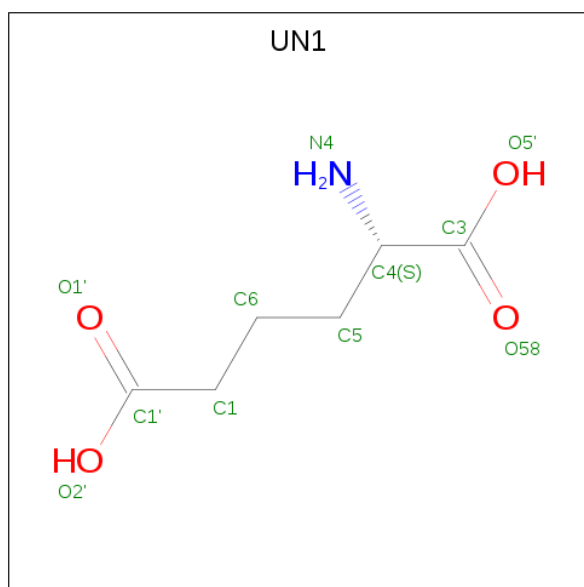
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P49419
A	0	HIS	-	expression tag	UNP P49419
A	399	ASP	GLU	engineered mutation	UNP P49419
B	-1	GLY	-	expression tag	UNP P49419
B	0	HIS	-	expression tag	UNP P49419
B	399	ASP	GLU	engineered mutation	UNP P49419
C	-1	GLY	-	expression tag	UNP P49419
C	0	HIS	-	expression tag	UNP P49419
C	399	ASP	GLU	engineered mutation	UNP P49419
D	-1	GLY	-	expression tag	UNP P49419
D	0	HIS	-	expression tag	UNP P49419
D	399	ASP	GLU	engineered mutation	UNP P49419
E	-1	GLY	-	expression tag	UNP P49419

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	expression tag	UNP P49419
E	399	ASP	GLU	engineered mutation	UNP P49419
F	-1	GLY	-	expression tag	UNP P49419
F	0	HIS	-	expression tag	UNP P49419
F	399	ASP	GLU	engineered mutation	UNP P49419
G	-1	GLY	-	expression tag	UNP P49419
G	0	HIS	-	expression tag	UNP P49419
G	399	ASP	GLU	engineered mutation	UNP P49419
H	-1	GLY	-	expression tag	UNP P49419
H	0	HIS	-	expression tag	UNP P49419
H	399	ASP	GLU	engineered mutation	UNP P49419

- Molecule 2 is 2-AMINOHEXANEDIOIC ACID (three-letter code: UN1) (formula: C₆H₁₁NO₄) (labeled as "Ligand of Interest" by author).



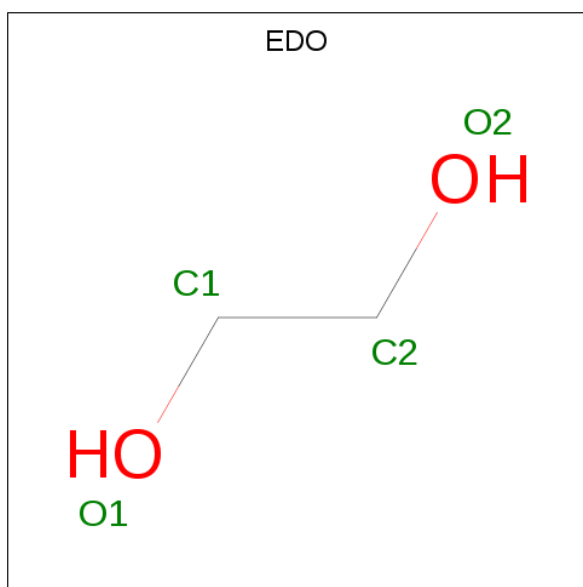
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	6	1	4		
2	B	1	Total	C	N	O	0	0
			11	6	1	4		
2	C	1	Total	C	N	O	0	0
			11	6	1	4		
2	D	1	Total	C	N	O	0	0
			11	6	1	4		
2	E	1	Total	C	N	O	0	0
			11	6	1	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	F	1	Total	C	N	O	0	0
			11	6	1	4		
2	G	1	Total	C	N	O	0	0
			11	6	1	4		
2	H	1	Total	C	N	O	0	0
			11	6	1	4		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	G	1	Total C O 4 2 2	0	0
3	G	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0

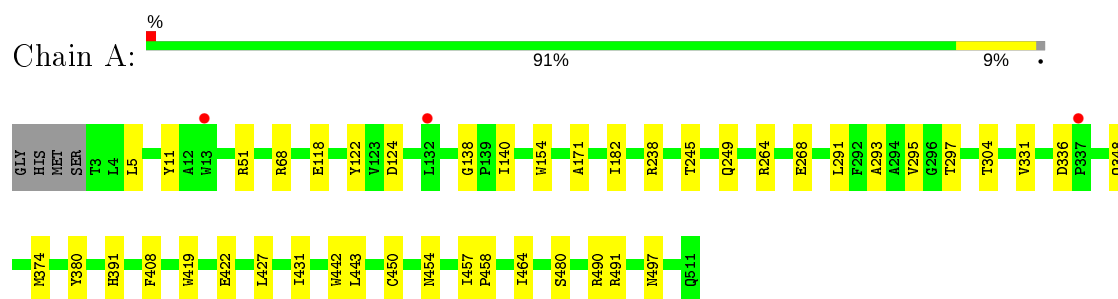
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	253	Total O 255 255	0	2
4	B	242	Total O 246 246	0	4
4	C	240	Total O 240 240	0	0
4	D	223	Total O 224 224	0	1
4	E	227	Total O 227 227	0	0
4	F	233	Total O 235 235	0	2
4	G	239	Total O 240 240	0	1
4	H	268	Total O 268 268	0	0

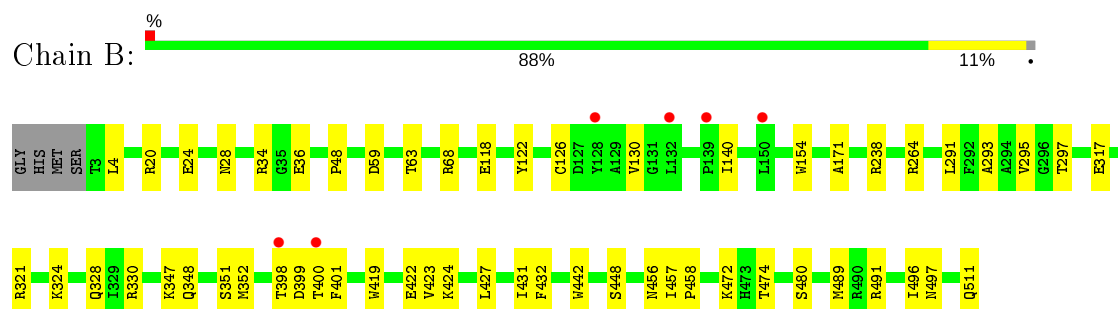
3 Residue-property plots [i](#)

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.

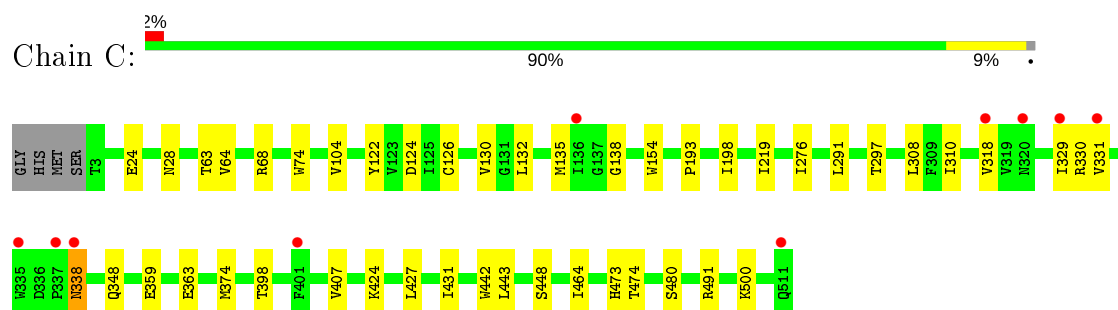
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



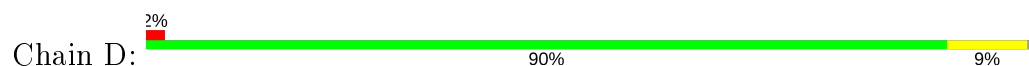
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase

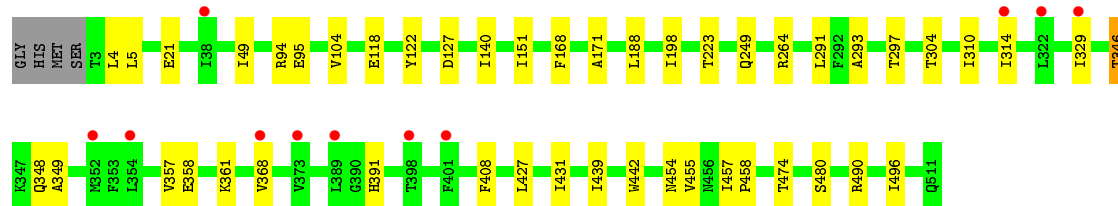


- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase

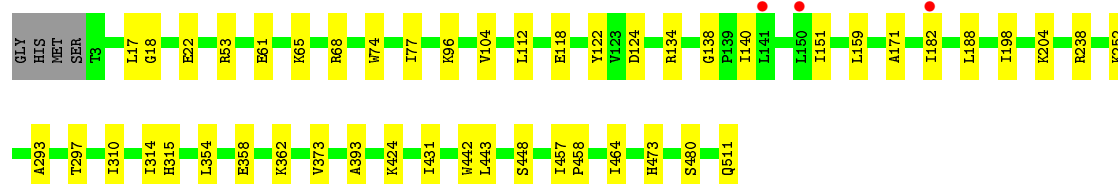
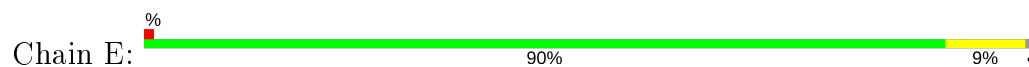


- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase

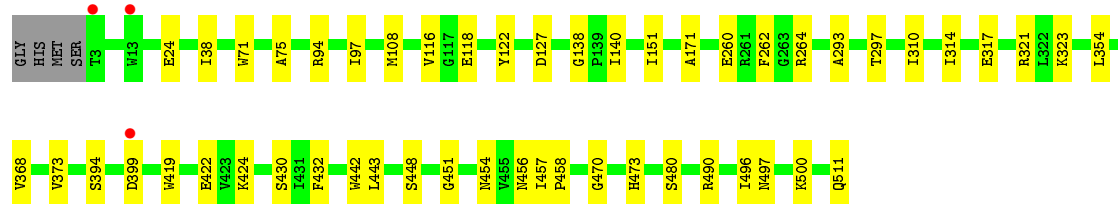
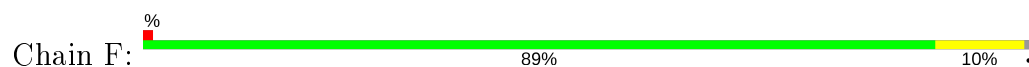




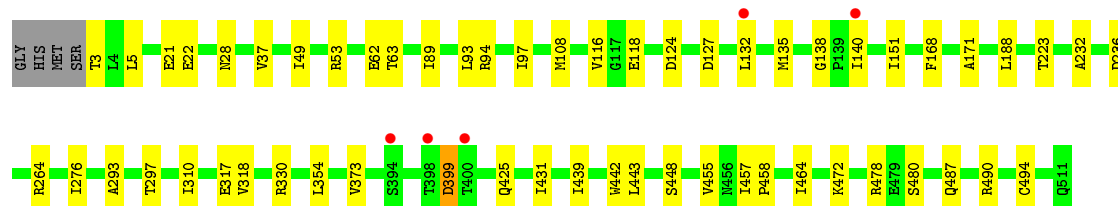
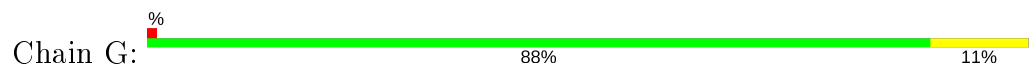
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



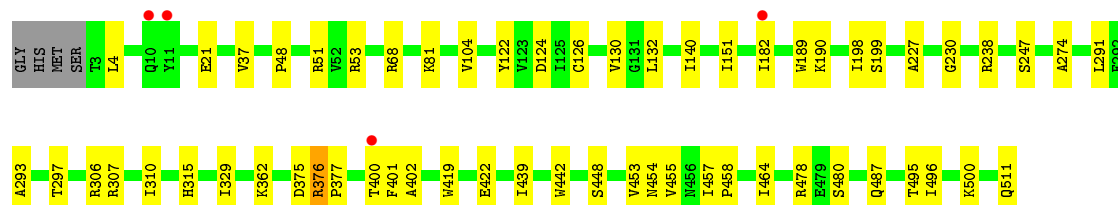
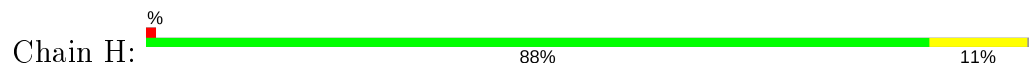
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.11Å 160.14Å 157.87Å 90.00° 95.07° 90.00°	Depositor
Resolution (Å)	49.03 – 1.75 49.03 – 1.75	Depositor EDS
% Data completeness (in resolution range)	94.8 (49.03-1.75) 98.6 (49.03-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.75Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.193 , 0.241 0.192 , 0.241	Depositor DCC
R_{free} test set	18899 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	33043	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.57% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: UN1, EDO

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	0.38	0/3949	0.53	0/5361
1	B	0.37	0/3949	0.55	0/5357
1	C	0.38	0/3949	0.54	0/5357
1	D	0.35	0/3947	0.52	0/5355
1	E	0.37	0/3967	0.54	0/5378
1	F	0.37	0/3961	0.53	0/5370
1	G	0.36	0/3957	0.53	0/5367
1	H	0.36	0/3953	0.53	0/5361
All	All	0.37	0/31632	0.53	0/42906

There are no bond length outliers.

There are no bond angle outliers.

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	3860	0	3827	27	0
1	B	3866	0	3850	30	0
1	C	3866	0	3850	24	0
1	D	3864	0	3843	25	0
1	E	3881	0	3883	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3878	0	3878	27	0
1	G	3871	0	3868	33	0
1	H	3870	0	3859	27	0
2	A	11	0	9	0	0
2	B	11	0	9	0	0
2	C	11	0	9	0	0
2	D	11	0	9	1	0
2	E	11	0	9	0	0
2	F	11	0	9	0	0
2	G	11	0	9	1	0
2	H	11	0	9	0	0
3	A	4	0	6	0	0
3	B	12	0	18	1	0
3	C	12	0	18	0	0
3	D	4	0	6	0	0
3	E	4	0	6	0	0
3	F	12	0	18	0	0
3	G	8	0	12	1	0
3	H	8	0	12	2	0
4	A	255	0	0	3	0
4	B	246	0	0	1	0
4	C	240	0	0	2	0
4	D	224	0	0	1	0
4	E	227	0	0	2	0
4	F	235	0	0	0	0
4	G	240	0	0	0	0
4	H	268	0	0	0	0
All	All	33043	0	31026	201	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 3.

All (201) 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:E:293:ALA:HB2	1:E:458:PRO:HB3	1.58	0.85
1:A:374:MET:HE2	1:A:380:TYR:HB3	1.64	0.80
1:F:317:GLU:OE2	1:F:321:ARG:NH1	2.22	0.72
1:E:443:LEU:HD11	1:F:496:ILE:HD11	1.71	0.72
1:G:118:GLU:HG3	1:G:171:ALA:HB2	1.72	0.70
1:G:478:ARG:O	1:G:487:GLN:NE2	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:TRP:HA	1:E:77:ILE:HD12	1.76	0.68
1:D:346:THR:HG23	1:D:349:ALA:H	1.59	0.68
1:F:118:GLU:HG3	1:F:171:ALA:HB2	1.76	0.66
1:D:264:ARG:HH22	1:D:490:ARG:HA	1.64	0.63
1:H:68:ARG:HD2	1:H:238:ARG:HB3	1.82	0.62
1:G:293:ALA:HB2	1:G:458:PRO:HB3	1.82	0.62
1:H:293:ALA:HB2	1:H:458:PRO:HB3	1.81	0.62
1:E:140:ILE:HD13	1:H:151:ILE:HG23	1.81	0.62
1:B:293:ALA:HB2	1:B:458:PRO:HB3	1.79	0.61
1:A:118:GLU:HG3	1:A:171:ALA:HB2	1.82	0.61
1:B:20:ARG:NH2	1:B:24:GLU:OE2	2.35	0.60
1:A:68:ARG:HD2	1:A:238:ARG:HB3	1.84	0.59
1:G:37:VAL:HG22	1:G:53:ARG:HG2	1.85	0.59
1:G:399:ASP:N	1:G:399:ASP:OD2	2.35	0.58
1:G:5:LEU:HB2	1:G:49:ILE:HA	1.86	0.58
1:D:94:ARG:NH2	1:D:127:ASP:OD1	2.32	0.58
1:D:293:ALA:HB2	1:D:458:PRO:HB3	1.86	0.58
1:G:22:GLU:HG2	1:G:53:ARG:HE	1.70	0.56
1:E:151:ILE:HD11	1:F:443:LEU:HD13	1.85	0.56
1:D:291:LEU:HD11	1:D:329:ILE:HD11	1.86	0.56
1:F:138:GLY:O	1:G:140[A]:ILE:HG22	2.04	0.56
1:B:28:ASN:HB3	1:B:63:THR:HG23	1.89	0.55
1:D:357:VAL:HG12	1:D:361:LYS:HE2	1.87	0.55
1:A:331:VAL:HG11	1:A:374:MET:CE	2.37	0.55
1:D:118:GLU:HG3	1:D:171:ALA:HB2	1.89	0.55
1:E:68:ARG:NH1	1:E:238:ARG:O	2.40	0.55
4:A:912:HOH:O	1:F:38:ILE:HD11	2.07	0.54
1:D:249:GLN:HB2	4:D:896:HOH:O	2.07	0.54
1:F:451:GLY:HA3	1:F:470:GLY:HA2	1.89	0.54
1:D:439:ILE:HG23	1:D:455:VAL:HG21	1.90	0.54
1:G:168:PHE:CE1	2:G:601:UN1:H6C2	2.43	0.53
1:E:159:LEU:HD11	1:E:188:LEU:HB2	1.89	0.53
1:B:427:LEU:HD13	1:B:474:THR:HG21	1.91	0.53
1:F:354:LEU:HD21	1:F:373:VAL:HG23	1.91	0.53
1:B:291:LEU:O	1:B:295:VAL:HG22	2.08	0.53
1:B:442:TRP:CH2	1:B:448:SER:HB2	2.43	0.53
1:E:431:ILE:HG23	1:E:442:TRP:CE2	2.43	0.52
1:A:419:TRP:O	1:A:422:GLU:HG2	2.10	0.52
1:H:230:GLY:HA3	3:H:603:EDO:H22	1.90	0.52
1:D:104:VAL:HA	1:D:198:ILE:HD11	1.90	0.52
1:E:118:GLU:HG3	1:E:171:ALA:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:28:ASN:HB3	1:G:63:THR:HG23	1.90	0.52
1:G:124:ASP:HB3	1:G:464:ILE:HG12	1.90	0.52
1:F:424:LYS:HD3	1:F:473:HIS:CE1	2.45	0.52
1:F:293:ALA:HB2	1:F:458:PRO:HB3	1.91	0.51
1:B:264:ARG:NH2	1:B:489:MET:O	2.43	0.51
1:E:354:LEU:HD21	1:E:373:VAL:HG23	1.93	0.51
1:B:317:GLU:OE2	1:B:321:ARG:NH1	2.43	0.51
1:G:443:LEU:HD11	1:H:496:ILE:HD11	1.92	0.51
1:C:331:VAL:HG11	1:C:374:MET:SD	2.50	0.51
1:G:264:ARG:HH22	1:G:490:ARG:HA	1.76	0.51
1:B:330:ARG:NH2	4:B:708:HOH:O	2.44	0.50
1:A:138:GLY:O	1:D:140:ILE:HG22	2.10	0.50
1:C:443:LEU:HD11	1:D:496:ILE:HD11	1.93	0.50
1:E:315:HIS:NE2	4:E:706:HOH:O	2.35	0.50
1:H:376:ARG:HG2	1:H:377:PRO:HD2	1.91	0.50
1:B:351:SER:HB2	1:E:354:LEU:HD12	1.92	0.50
1:E:442:TRP:CH2	1:E:448:SER:HB2	2.46	0.50
1:C:330:ARG:NH2	1:C:338:ASN:O	2.45	0.50
1:E:124:ASP:HB3	1:E:464:ILE:HG12	1.93	0.50
1:G:457:ILE:HD11	1:H:495:THR:HB	1.93	0.50
1:G:89:ILE:O	1:G:93:LEU:HD23	2.12	0.50
1:A:457:ILE:HG23	1:B:497:ASN:HB2	1.93	0.50
1:G:132:LEU:HD23	1:G:135:MET:CE	2.42	0.50
1:E:138:GLY:O	1:H:140:ILE:HG22	2.12	0.50
1:G:276:ILE:HB	1:G:431:ILE:HG22	1.94	0.49
1:G:442:TRP:CH2	1:G:448:SER:HB2	2.47	0.49
1:H:419:TRP:O	1:H:422:GLU:HG2	2.12	0.49
1:D:4:LEU:HD12	1:D:21:GLU:HG2	1.95	0.49
1:C:424:LYS:HD3	1:C:473:HIS:CE1	2.47	0.49
1:B:431:ILE:HG23	1:B:442:TRP:CE2	2.48	0.49
1:B:419:TRP:O	1:B:422:GLU:HG2	2.12	0.49
1:C:132:LEU:O	1:C:135:MET:N	2.36	0.49
1:E:17:LEU:HD23	1:E:96:LYS:HD2	1.95	0.48
1:H:291:LEU:HD11	1:H:329:ILE:HD11	1.96	0.48
1:A:374:MET:CE	1:A:380:TYR:HB3	2.41	0.48
1:C:310:ILE:HD12	1:C:407:VAL:HG13	1.95	0.48
1:F:264:ARG:HH12	1:F:490:ARG:HA	1.79	0.48
1:F:419:TRP:O	1:F:422:GLU:HG2	2.13	0.48
1:G:354:LEU:HD21	1:G:373:VAL:HG23	1.96	0.48
1:C:310:ILE:HG21	1:C:318:VAL:HG21	1.96	0.48
1:H:4:LEU:HD23	1:H:48:PRO:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:358:GLU:O	1:E:362:LYS:HG3	2.14	0.47
1:C:154:TRP:CH2	1:C:491:ARG:HB2	2.49	0.47
1:A:331:VAL:HG11	1:A:374:MET:HE3	1.97	0.47
1:A:140:ILE:HD13	1:D:151:ILE:HG23	1.96	0.47
1:H:454:ASN:HB3	1:H:457:ILE:HG13	1.96	0.47
1:A:293:ALA:HB2	1:A:458:PRO:HB3	1.97	0.47
1:B:118:GLU:HG3	1:B:171:ALA:HB2	1.96	0.47
1:D:391:HIS:HA	1:D:408:PHE:CE1	2.49	0.47
1:H:126:CYS:O	1:H:130:VAL:HG23	2.13	0.47
1:H:124:ASP:HB3	1:H:464:ILE:HG12	1.96	0.47
1:H:274:ALA:HA	1:H:307:ARG:O	2.15	0.47
1:A:154:TRP:CH2	1:A:491:ARG:HB2	2.50	0.47
1:C:104:VAL:HA	1:C:198:ILE:HD11	1.96	0.47
1:G:135:MET:HE2	1:G:135:MET:HB2	1.31	0.46
1:E:61:GLU:HG3	1:E:65:LYS:HE3	1.98	0.46
1:H:37:VAL:HG22	1:H:53:ARG:HG2	1.97	0.46
1:A:124:ASP:HB3	1:A:464:ILE:HG12	1.97	0.46
1:A:249:GLN:HB2	4:A:897:HOH:O	2.16	0.46
1:F:432:PHE:CD2	1:F:456:ASN:HA	2.50	0.46
1:B:34:ARG:HD3	1:B:59:ASP:OD1	2.16	0.45
1:H:310:ILE:HG21	1:H:315:HIS:HA	1.98	0.45
1:A:497:ASN:HB2	1:B:457:ILE:HG23	1.97	0.45
1:B:140:ILE:HG22	1:C:138:GLY:O	2.15	0.45
1:E:424:LYS:HD3	1:E:473:HIS:CE1	2.52	0.45
1:G:140[A]:ILE:HG13	1:G:151:ILE:HG12	1.98	0.45
1:H:439:ILE:HG23	1:H:455:VAL:HG21	1.99	0.45
1:A:5:LEU:HD22	1:A:11:TYR:CE2	2.51	0.45
1:G:62:GLU:HG3	3:G:603:EDO:H11	1.98	0.45
1:B:68:ARG:HD2	1:B:238:ARG:HB3	1.99	0.45
1:D:431:ILE:HG12	1:D:442:TRP:CD1	2.51	0.45
1:E:22:GLU:HG2	1:E:53:ARG:HE	1.82	0.45
1:F:310:ILE:HG22	1:F:314:ILE:HG13	1.98	0.45
1:H:442:TRP:CH2	1:H:448:SER:HB2	2.52	0.45
1:F:94:ARG:NH1	1:F:127:ASP:OD2	2.50	0.45
1:G:425:GLN:O	1:G:472:LYS:HE3	2.16	0.45
1:C:500:LYS:HE3	4:C:753:HOH:O	2.17	0.44
1:E:140:ILE:HD11	1:H:140:ILE:HB	1.98	0.44
1:B:36:GLU:OE2	1:E:393:ALA:HA	2.16	0.44
1:D:427:LEU:HD13	1:D:474:THR:HG21	2.00	0.44
1:A:291:LEU:O	1:A:295:VAL:HG22	2.17	0.44
1:D:310:ILE:HG22	1:D:314:ILE:HG13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:104:VAL:HA	1:E:198:ILE:HD11	1.99	0.44
1:G:439:ILE:HG23	1:G:455:VAL:HG21	2.00	0.44
1:H:478:ARG:O	1:H:487:GLN:NE2	2.39	0.44
1:B:348:GLN:HG2	1:B:352:MET:HE2	2.00	0.44
1:D:454:ASN:HB3	1:D:457:ILE:HG13	1.99	0.44
1:H:400:THR:HG22	1:H:402:ALA:H	1.82	0.43
1:A:431:ILE:HG23	1:A:442:TRP:CE2	2.53	0.43
1:D:168:PHE:CE1	2:D:601:UN1:H6C2	2.53	0.43
1:F:500:LYS:HD3	1:F:500:LYS:HA	1.60	0.43
1:F:71:TRP:CE2	1:F:75:ALA:HB2	2.53	0.43
1:H:104:VAL:HA	1:H:198:ILE:HD11	2.01	0.43
1:G:97:ILE:HG12	1:G:116:VAL:HG13	2.00	0.43
1:A:427:LEU:O	1:A:450:CYS:HB3	2.19	0.43
1:E:18:GLY:O	1:E:204:LYS:HE2	2.18	0.43
1:D:188:LEU:HD11	1:D:223:THR:OG1	2.19	0.43
1:F:442:TRP:CH2	1:F:448:SER:HB2	2.54	0.43
1:A:331:VAL:HG11	1:A:374:MET:HE1	1.99	0.43
1:E:310:ILE:HG22	1:E:314:ILE:HG13	1.99	0.43
1:F:97:ILE:HG12	1:F:116:VAL:HG13	2.00	0.43
1:C:308:LEU:HD23	1:C:407:VAL:HG22	2.01	0.43
1:D:358:GLU:HA	1:D:361:LYS:HE3	2.01	0.43
1:E:134:ARG:HG3	4:E:747:HOH:O	2.19	0.43
1:E:252:LYS:HB2	1:F:262:PHE:CZ	2.54	0.43
1:G:132:LEU:HD23	1:G:135:MET:HE1	2.00	0.43
1:B:4:LEU:HD23	1:B:48:PRO:HB2	2.00	0.43
1:E:68:ARG:HA	1:E:68:ARG:HD3	1.76	0.43
1:B:324:LYS:O	1:B:328:GLN:HG2	2.19	0.42
1:C:124:ASP:HB3	1:C:464:ILE:HG12	2.00	0.42
1:F:323:LYS:HD2	1:F:368:VAL:O	2.19	0.42
1:A:443:LEU:HD11	1:B:496:ILE:HD11	2.01	0.42
1:B:423:VAL:O	1:B:472:LYS:NZ	2.53	0.42
1:G:310:ILE:HD13	1:G:318:VAL:HB	2.00	0.42
1:A:454:ASN:HB3	1:A:457:ILE:HG13	2.01	0.42
1:C:310:ILE:HD13	1:C:318:VAL:CG2	2.50	0.42
1:C:359:GLU:HG3	1:C:363:GLU:OE1	2.20	0.42
1:G:431:ILE:HG23	1:G:442:TRP:CZ2	2.54	0.42
1:H:227:ALA:HA	3:H:603:EDO:H21	2.02	0.42
1:D:5:LEU:HB2	1:D:49:ILE:HA	2.01	0.42
1:F:454:ASN:HB3	1:F:457:ILE:HG13	2.02	0.42
1:G:94:ARG:NH2	1:G:127:ASP:OD1	2.47	0.42
1:D:431:ILE:HG23	1:D:442:TRP:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:494:CYS:HA	1:H:453:VAL:O	2.20	0.41
1:C:64:VAL:O	1:C:68:ARG:HG2	2.20	0.41
1:C:74:TRP:CE3	1:C:219:ILE:HD13	2.55	0.41
1:A:51:ARG:HD3	4:A:871:HOH:O	2.21	0.41
1:B:126:CYS:O	1:B:130:VAL:HG23	2.20	0.41
1:B:424:LYS:HE2	1:B:424:LYS:HA	2.01	0.41
1:C:126:CYS:O	1:C:130:VAL:HG23	2.20	0.41
1:B:432:PHE:CD2	1:B:456:ASN:HA	2.54	0.41
1:C:28:ASN:HB3	1:C:63:THR:HG23	2.02	0.41
1:E:443:LEU:HD13	1:F:151:ILE:HD11	2.03	0.41
1:A:293:ALA:O	1:A:304:THR:HG23	2.21	0.41
1:A:264:ARG:HH22	1:A:490:ARG:HA	1.86	0.41
1:A:391:HIS:HA	1:A:408:PHE:CE1	2.56	0.41
1:E:252:LYS:HE2	1:F:260:GLU:O	2.21	0.41
1:G:188:LEU:HD11	1:G:223:THR:HG23	2.03	0.41
1:A:245:THR:HA	1:A:268:GLU:O	2.21	0.41
1:C:276:ILE:HB	1:C:431:ILE:HG22	2.02	0.41
1:H:21:GLU:HG2	1:H:51:ARG:NH2	2.36	0.41
1:C:442:TRP:CH2	1:C:448:SER:HB2	2.56	0.40
1:C:427:LEU:HD13	1:C:474:THR:HG21	2.02	0.40
1:B:347:LYS:HD2	1:E:358:GLU:CG	2.52	0.40
1:C:291:LEU:HD21	1:C:329:ILE:HD11	2.02	0.40
1:F:430:SER:HA	1:F:454:ASN:O	2.22	0.40
1:H:189:TRP:CH2	1:H:199:SER:HA	2.57	0.40
1:B:401:PHE:HB2	3:B:603:EDO:H12	2.04	0.40
1:D:293:ALA:O	1:D:304:THR:HG23	2.21	0.40
1:F:140:ILE:HG22	1:G:138:GLY:O	2.21	0.40
1:B:154:TRP:CH2	1:B:491:ARG:HB2	2.56	0.40
1:C:193:PRO:HD2	4:C:913:HOH:O	2.21	0.40
1:E:457:ILE:HG23	1:F:497:ASN:HB2	2.04	0.40
1:G:232:ALA:O	1:G:236:ASP:HB2	2.22	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	510/513 (99%)	494 (97%)	15 (3%)	1 (0%)	47	29
1	B	509/513 (99%)	494 (97%)	14 (3%)	1 (0%)	47	29
1	C	508/513 (99%)	486 (96%)	21 (4%)	1 (0%)	47	29
1	D	508/513 (99%)	494 (97%)	13 (3%)	1 (0%)	47	29
1	E	509/513 (99%)	490 (96%)	18 (4%)	1 (0%)	47	29
1	F	508/513 (99%)	487 (96%)	20 (4%)	1 (0%)	47	29
1	G	509/513 (99%)	490 (96%)	18 (4%)	1 (0%)	47	29
1	H	508/513 (99%)	489 (96%)	18 (4%)	1 (0%)	47	29
All	All	4069/4104 (99%)	3924 (96%)	137 (3%)	8 (0%)	47	29

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	480	SER
1	B	480	SER
1	C	480	SER
1	D	480	SER
1	F	480	SER
1	G	480	SER
1	H	480	SER
1	E	480	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	402/410 (98%)	397 (99%)	5 (1%)	71	56
1	B	403/410 (98%)	397 (98%)	6 (2%)	65	49
1	C	403/410 (98%)	397 (98%)	6 (2%)	65	49
1	D	403/410 (98%)	397 (98%)	6 (2%)	65	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	407/410 (99%)	402 (99%)	5 (1%)	71	56
1	F	406/410 (99%)	399 (98%)	7 (2%)	60	42
1	G	404/410 (98%)	397 (98%)	7 (2%)	60	42
1	H	404/410 (98%)	390 (96%)	14 (4%)	36	13
All	All	3232/3280 (98%)	3176 (98%)	56 (2%)	60	42

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	TYR
1	A	182	ILE
1	A	297	THR
1	A	336	ASP
1	A	348	GLN
1	B	122	TYR
1	B	297	THR
1	B	398	THR
1	B	399	ASP
1	B	400	THR
1	B	511	GLN
1	C	24	GLU
1	C	122	TYR
1	C	297	THR
1	C	338	ASN
1	C	348	GLN
1	C	398	THR
1	D	95	GLU
1	D	122	TYR
1	D	297	THR
1	D	346	THR
1	D	348	GLN
1	D	368	VAL
1	E	112	LEU
1	E	122	TYR
1	E	182	ILE
1	E	297	THR
1	E	511	GLN
1	F	24	GLU
1	F	108	MET
1	F	122	TYR
1	F	297	THR

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Mol	Chain	Res	Type
1	F	394	SER
1	F	399	ASP
1	F	511	GLN
1	G	3	THR
1	G	21	GLU
1	G	108	MET
1	G	297	THR
1	G	317	GLU
1	G	330	ARG
1	G	399	ASP
1	H	81	LYS
1	H	122	TYR
1	H	132	LEU
1	H	182	ILE
1	H	190	LYS
1	H	247	SER
1	H	297	THR
1	H	306	ARG
1	H	362	LYS
1	H	375	ASP
1	H	376	ARG
1	H	401	PHE
1	H	500	LYS
1	H	511	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	473	HIS
1	F	473	HIS

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 ⓘ

24 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	UN1	F	601	-	3,10,10	0.39	0	2,12,12	1.02	0
2	UN1	H	601	-	3,10,10	0.39	0	2,12,12	1.06	0
2	UN1	G	601	-	3,10,10	0.51	0	2,12,12	0.69	0
3	EDO	H	603	-	3,3,3	0.46	0	2,2,2	0.14	0
2	UN1	B	601	-	3,10,10	0.46	0	2,12,12	1.95	1 (50%)
3	EDO	F	602	-	3,3,3	0.47	0	2,2,2	0.23	0
2	UN1	D	601	-	3,10,10	0.36	0	2,12,12	0.77	0
2	UN1	C	601	-	3,10,10	0.50	0	2,12,12	0.69	0
2	UN1	E	601	-	3,10,10	0.48	0	2,12,12	0.47	0
3	EDO	G	603	-	3,3,3	0.49	0	2,2,2	0.33	0
3	EDO	H	602	-	3,3,3	0.47	0	2,2,2	0.37	0
3	EDO	C	604	-	3,3,3	0.46	0	2,2,2	0.32	0
3	EDO	B	603	-	3,3,3	0.47	0	2,2,2	0.24	0
3	EDO	D	602	-	3,3,3	0.46	0	2,2,2	0.35	0
3	EDO	G	602	-	3,3,3	0.48	0	2,2,2	0.20	0
3	EDO	E	602	-	3,3,3	0.46	0	2,2,2	0.43	0
3	EDO	C	602	-	3,3,3	0.50	0	2,2,2	0.28	0
3	EDO	A	602	-	3,3,3	0.48	0	2,2,2	0.37	0
3	EDO	B	604	-	3,3,3	0.49	0	2,2,2	0.17	0
3	EDO	C	603	-	3,3,3	0.51	0	2,2,2	0.16	0
3	EDO	F	603	-	3,3,3	0.47	0	2,2,2	0.31	0
3	EDO	B	602	-	3,3,3	0.47	0	2,2,2	0.26	0
2	UN1	A	601	-	3,10,10	0.39	0	2,12,12	0.77	0
3	EDO	F	604	-	3,3,3	0.49	0	2,2,2	0.22	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
2	UN1	F	601	-	-	0/4/10/10	-
2	UN1	H	601	-	-	0/4/10/10	-
2	UN1	G	601	-	-	0/4/10/10	-
3	EDO	H	603	-	-	1/1/1/1	-
2	UN1	B	601	-	-	0/4/10/10	-
3	EDO	F	602	-	-	1/1/1/1	-
2	UN1	D	601	-	-	0/4/10/10	-
2	UN1	C	601	-	-	0/4/10/10	-
2	UN1	E	601	-	-	0/4/10/10	-
3	EDO	G	603	-	-	0/1/1/1	-
3	EDO	H	602	-	-	1/1/1/1	-
3	EDO	C	604	-	-	1/1/1/1	-
3	EDO	B	603	-	-	1/1/1/1	-
3	EDO	D	602	-	-	1/1/1/1	-
3	EDO	G	602	-	-	0/1/1/1	-
3	EDO	E	602	-	-	1/1/1/1	-
3	EDO	C	602	-	-	0/1/1/1	-
3	EDO	A	602	-	-	1/1/1/1	-
3	EDO	B	604	-	-	0/1/1/1	-
3	EDO	C	603	-	-	1/1/1/1	-
3	EDO	F	603	-	-	1/1/1/1	-
3	EDO	B	602	-	-	1/1/1/1	-
2	UN1	A	601	-	-	0/4/10/10	-
3	EDO	F	604	-	-	0/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	UN1	C5-C6-C1	-2.68	106.75	112.65

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	603	EDO	O1-C1-C2-O2
3	E	602	EDO	O1-C1-C2-O2
3	F	603	EDO	O1-C1-C2-O2
3	B	602	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	602	EDO	O1-C1-C2-O2
3	C	603	EDO	O1-C1-C2-O2
3	F	602	EDO	O1-C1-C2-O2
3	C	604	EDO	O1-C1-C2-O2
3	B	603	EDO	O1-C1-C2-O2
3	H	602	EDO	O1-C1-C2-O2
3	D	602	EDO	O1-C1-C2-O2

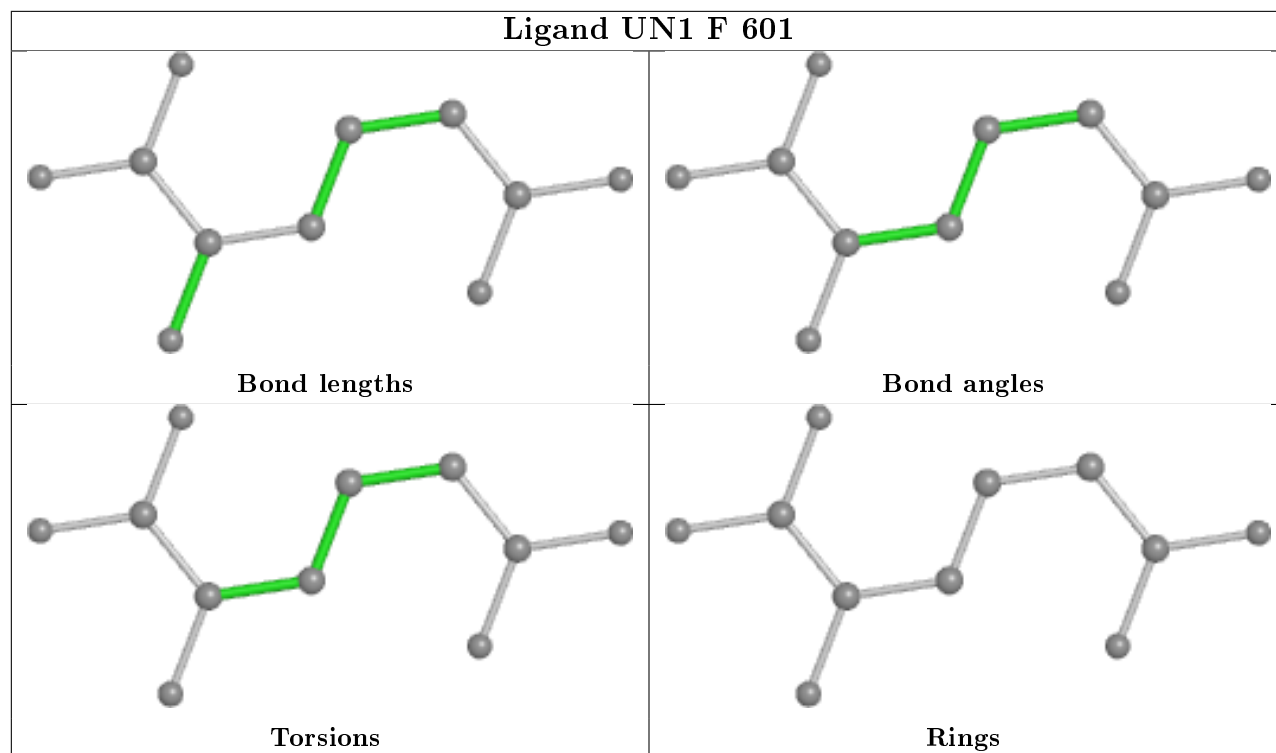
There are no ring outliers.

5 monomers are involved in 6 short contacts:

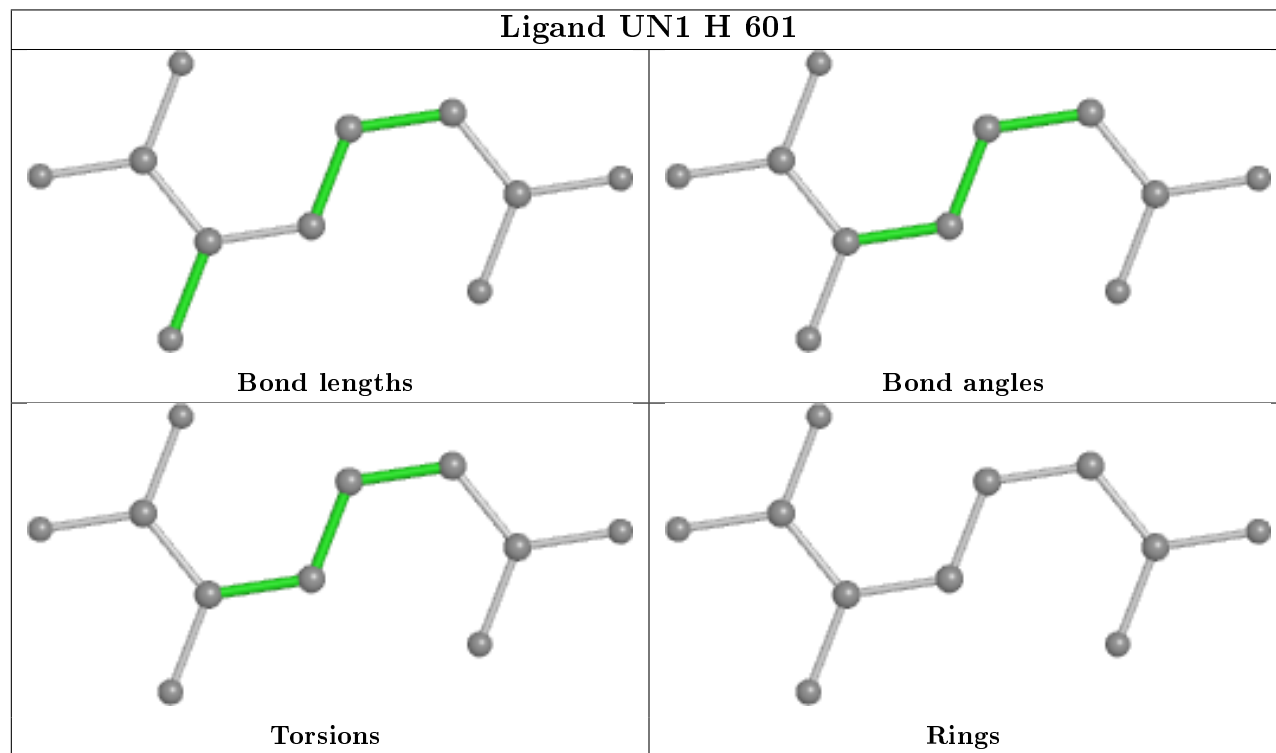
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	601	UN1	1	0
3	H	603	EDO	2	0
2	D	601	UN1	1	0
3	G	603	EDO	1	0
3	B	603	EDO	1	0

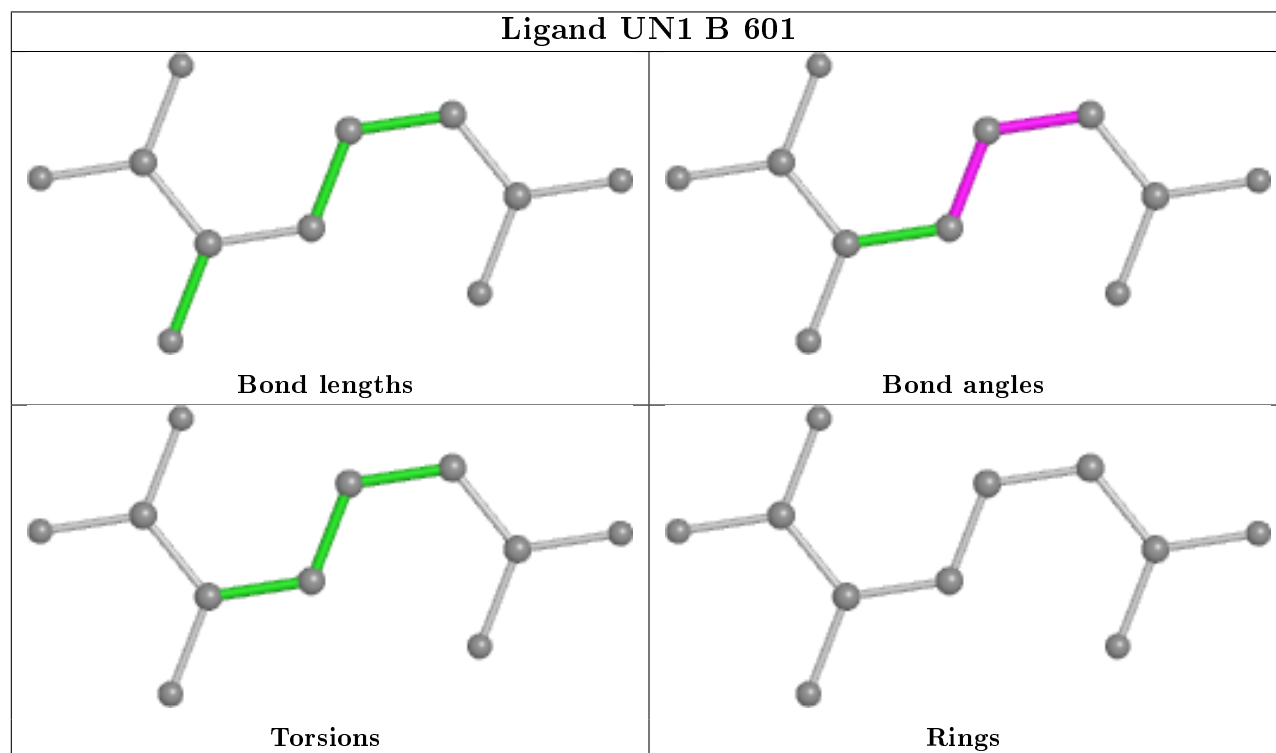
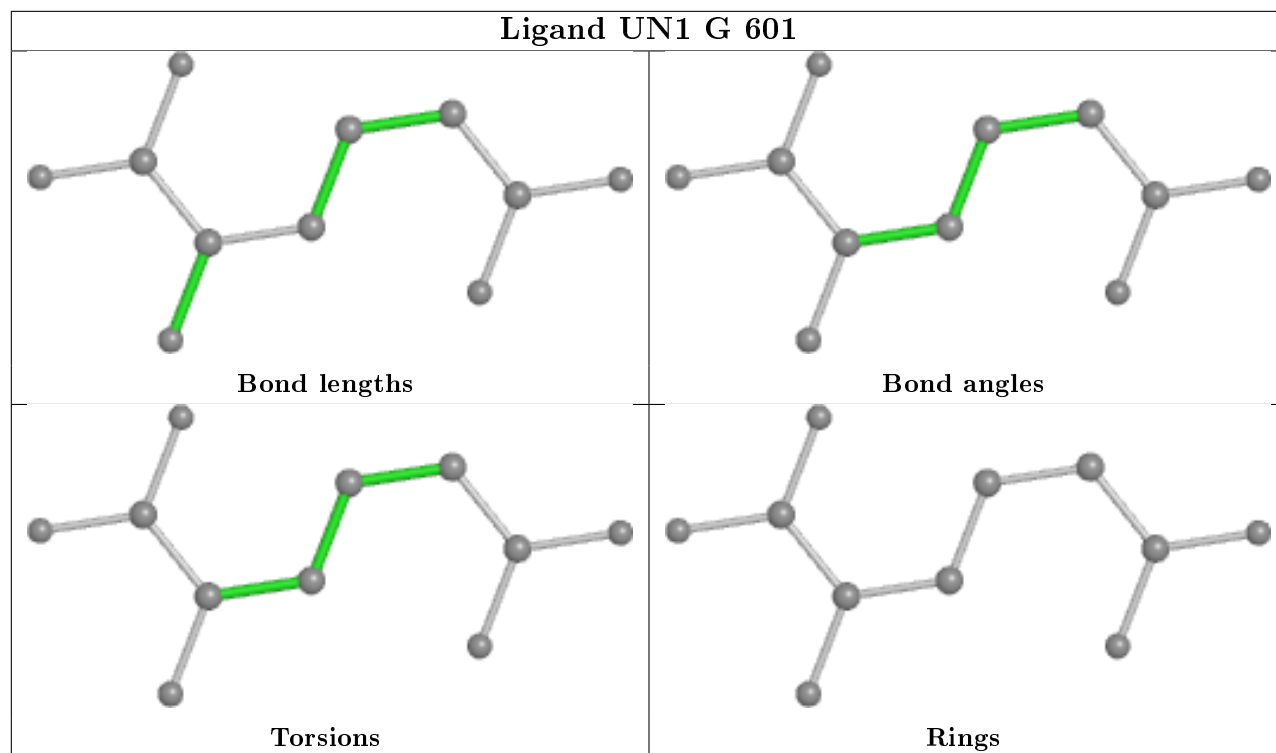
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

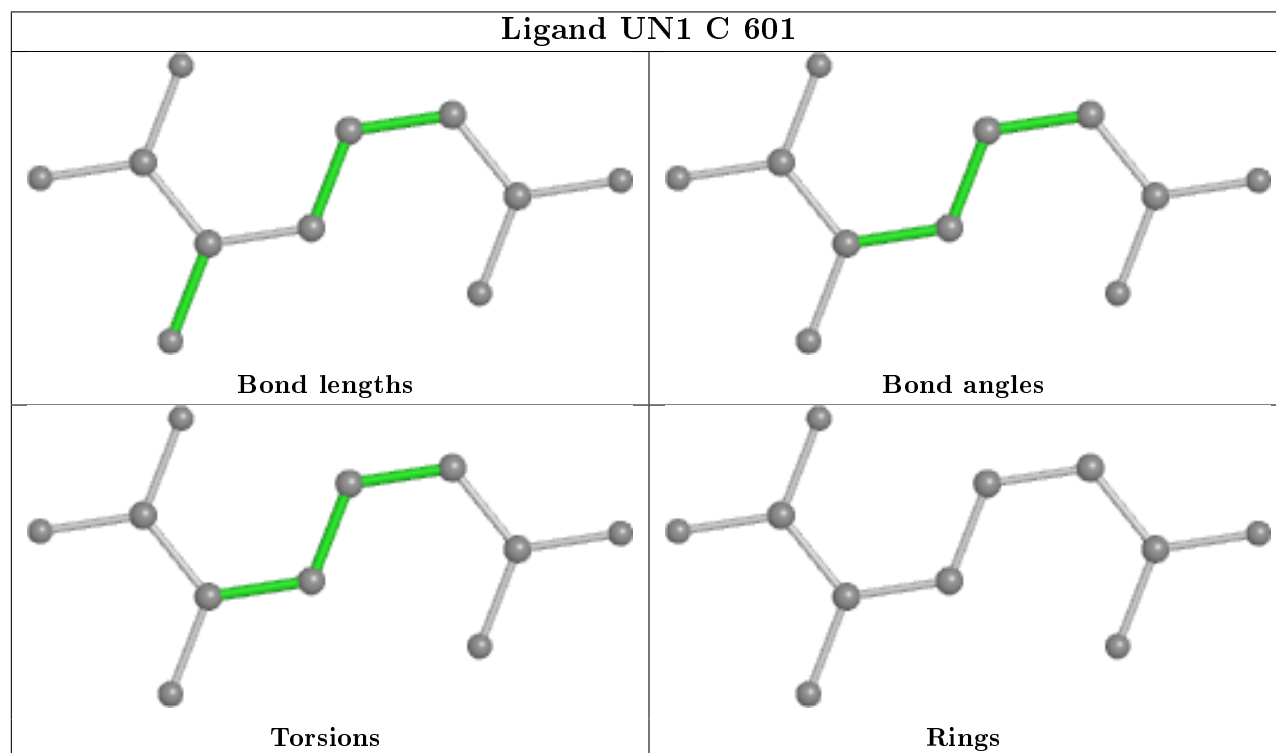
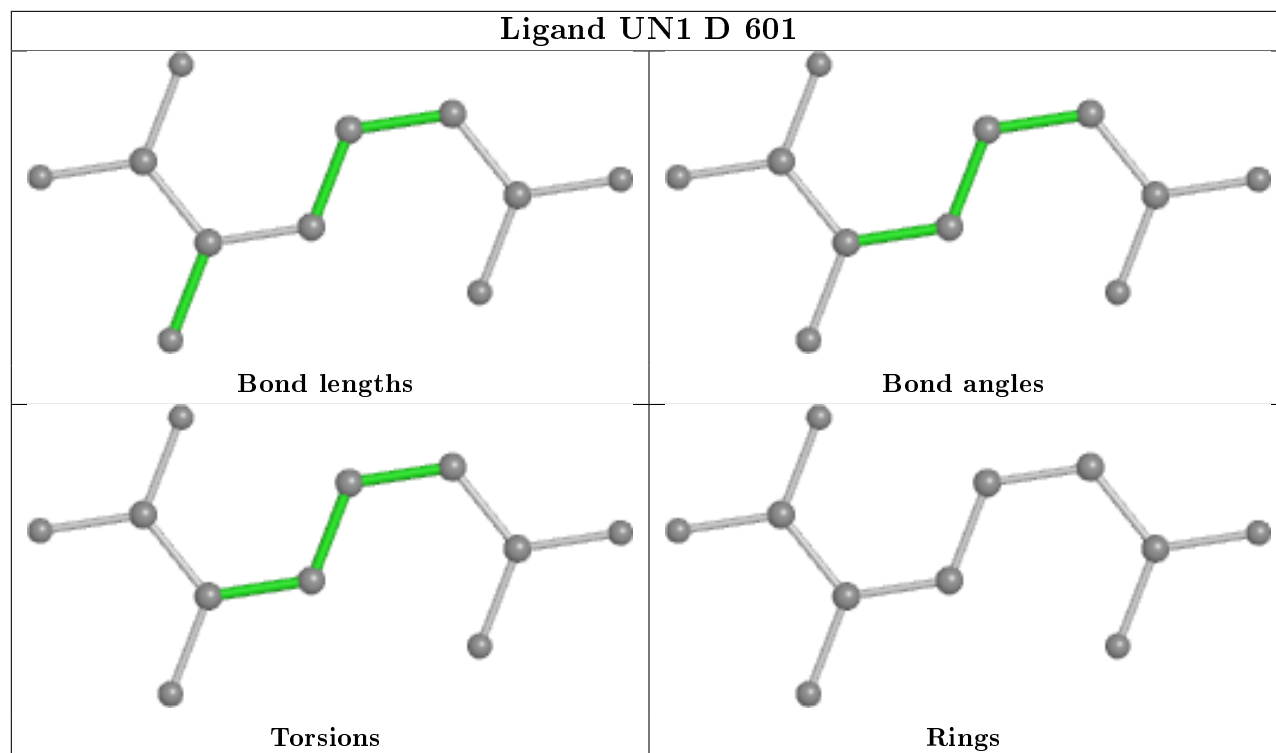
Ligand UN1 F 601

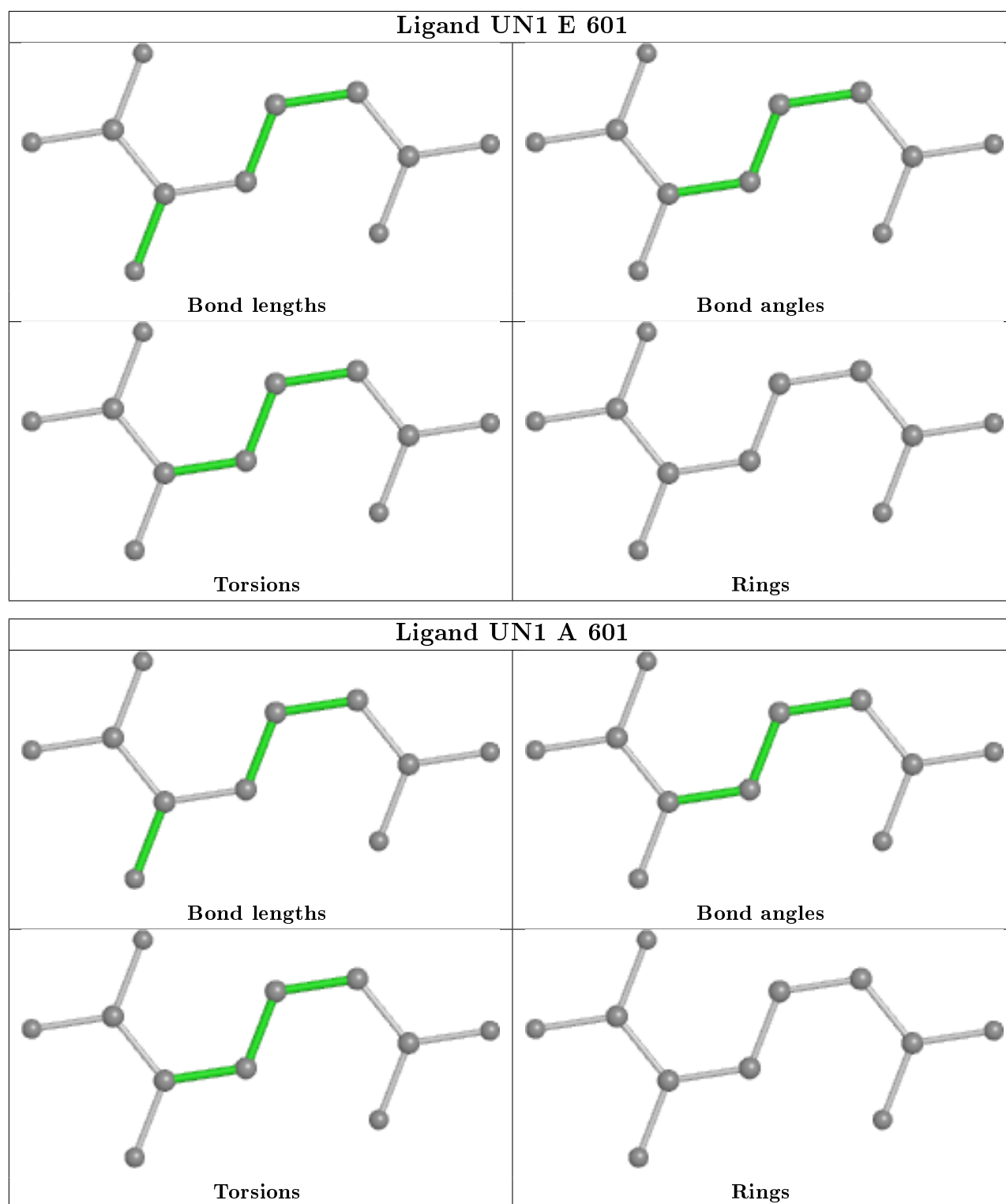


Ligand UN1 H 601









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	509/513 (99%)	0.02	3 (0%) 89 92	25, 35, 49, 70	0
1	B	509/513 (99%)	-0.01	6 (1%) 79 84	25, 35, 51, 84	0
1	C	509/513 (99%)	0.15	10 (1%) 65 72	24, 36, 61, 88	0
1	D	509/513 (99%)	0.12	11 (2%) 62 69	24, 37, 58, 88	0
1	E	509/513 (99%)	0.00	3 (0%) 89 92	26, 37, 55, 73	0
1	F	509/513 (99%)	0.05	3 (0%) 89 92	27, 37, 54, 98	0
1	G	509/513 (99%)	0.00	5 (0%) 82 87	29, 37, 53, 93	0
1	H	509/513 (99%)	0.01	4 (0%) 86 90	25, 34, 54, 89	0
All	All	4072/4104 (99%)	0.04	45 (1%) 80 86	24, 36, 55, 98	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	335	TRP	4.6
1	G	398	THR	3.7
1	G	400	THR	3.1
1	D	373	VAL	3.1
1	C	136	ILE	3.1
1	B	132	LEU	3.0
1	C	511	GLN	3.0
1	F	399	ASP	2.9
1	G	394	SER	2.9
1	B	398	THR	2.9
1	H	400	THR	2.9
1	G	140[A]	ILE	2.9
1	C	331	VAL	2.7
1	C	329	ILE	2.7
1	E	141	LEU	2.7
1	E	150	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	F	13	TRP	2.6
1	D	322	LEU	2.5
1	D	352	MET	2.5
1	A	13	TRP	2.4
1	C	337	PRO	2.3
1	D	389	LEU	2.3
1	H	11	TYR	2.3
1	H	10	GLN	2.3
1	C	401	PHE	2.3
1	C	318	VAL	2.3
1	D	398	THR	2.3
1	B	400	THR	2.2
1	H	182	ILE	2.2
1	D	329	ILE	2.2
1	C	320	ASN	2.2
1	A	132	LEU	2.2
1	D	368	VAL	2.2
1	D	354	LEU	2.2
1	D	314	ILE	2.2
1	D	401	PHE	2.1
1	D	38	ILE	2.1
1	B	128	TYR	2.1
1	C	338	ASN	2.1
1	A	337	PRO	2.1
1	B	150	LEU	2.1
1	E	182	ILE	2.0
1	B	139	PRO	2.0
1	G	132	LEU	2.0
1	F	3	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

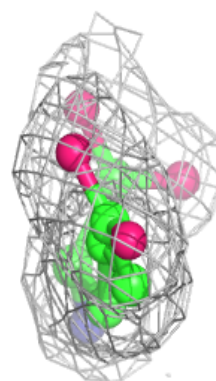
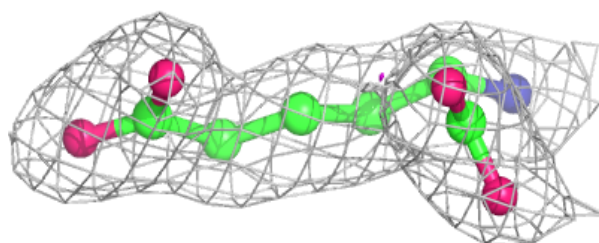
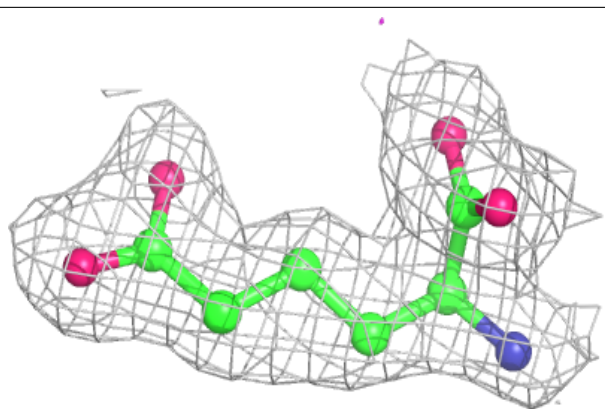
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	EDO	C	604	4/4	0.76	0.17	59,60,62,62	0
3	EDO	G	603	4/4	0.82	0.16	63,63,67,67	0
3	EDO	B	603	4/4	0.84	0.23	40,51,51,53	0
3	EDO	A	602	4/4	0.84	0.32	48,59,60,61	0
3	EDO	C	603	4/4	0.85	0.12	52,53,53,56	0
3	EDO	G	602	4/4	0.86	0.38	48,50,51,54	0
3	EDO	H	602	4/4	0.87	0.14	41,46,50,55	0
3	EDO	C	602	4/4	0.87	0.24	59,59,60,62	0
3	EDO	B	602	4/4	0.87	0.17	47,55,55,55	0
3	EDO	F	604	4/4	0.87	0.14	40,41,44,49	0
3	EDO	F	603	4/4	0.89	0.16	56,57,59,60	0
2	UN1	E	601	11/11	0.91	0.08	32,35,41,42	0
2	UN1	C	601	11/11	0.91	0.12	22,37,43,44	0
3	EDO	E	602	4/4	0.92	0.11	40,40,41,52	0
3	EDO	B	604	4/4	0.92	0.13	41,46,52,54	0
3	EDO	D	602	4/4	0.92	0.08	49,50,53,54	0
3	EDO	H	603	4/4	0.93	0.10	35,40,44,48	0
2	UN1	A	601	11/11	0.93	0.11	32,39,44,45	0
3	EDO	F	602	4/4	0.93	0.10	46,46,50,50	0
2	UN1	G	601	11/11	0.94	0.09	28,33,40,43	0
2	UN1	H	601	11/11	0.95	0.07	29,31,42,47	0
2	UN1	D	601	11/11	0.96	0.07	30,33,44,45	0
2	UN1	F	601	11/11	0.96	0.09	25,33,36,38	0
2	UN1	B	601	11/11	0.97	0.08	26,33,39,40	0

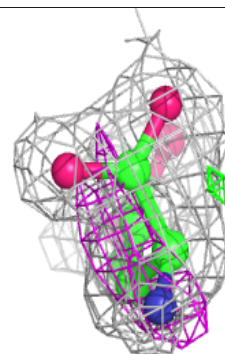
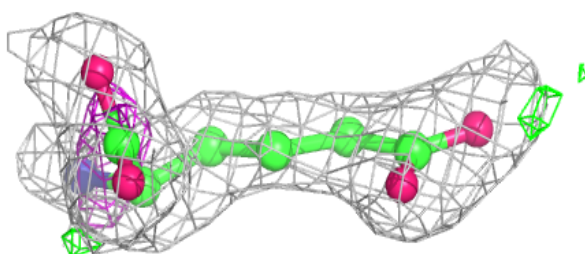
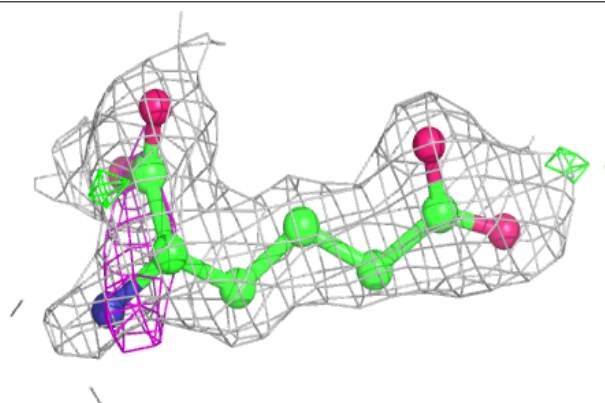
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around UN1 E 601:

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

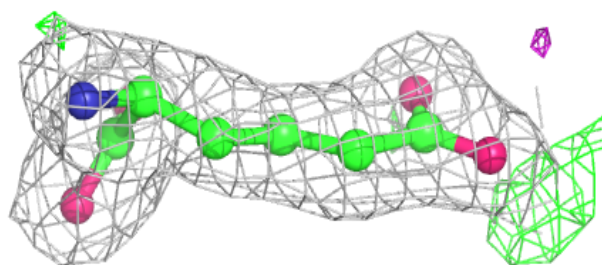
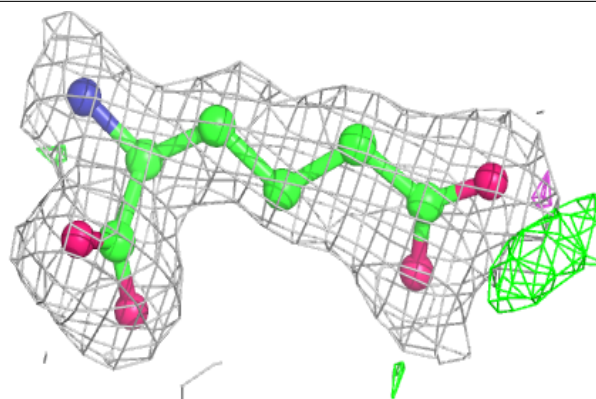
**Electron density around UN1 C 601:**

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

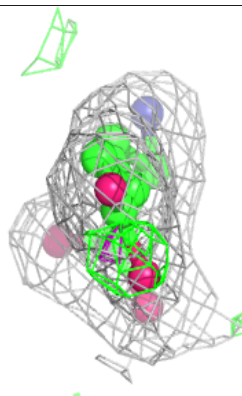
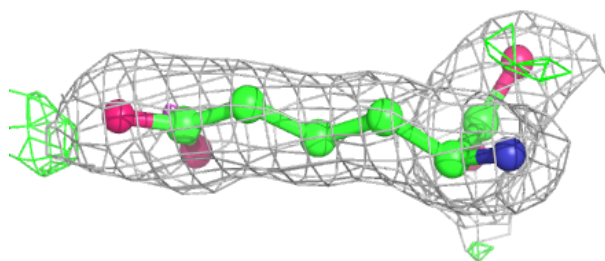
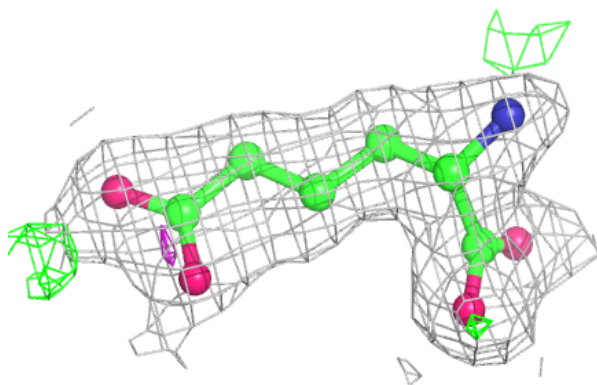


Electron density around UN1 A 601:

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

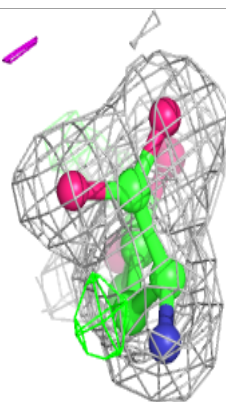
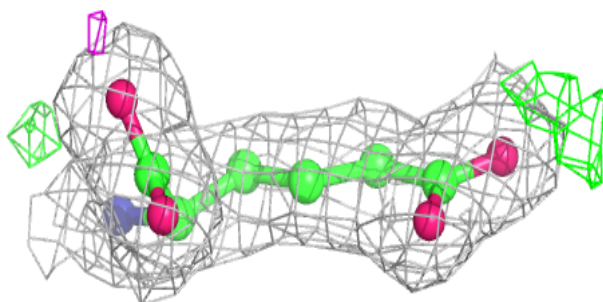
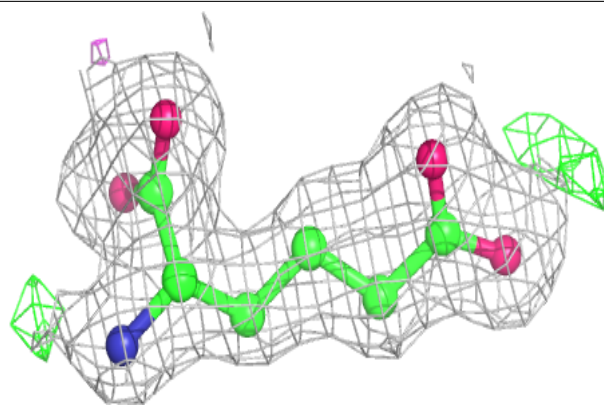
**Electron density around UN1 G 601:**

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

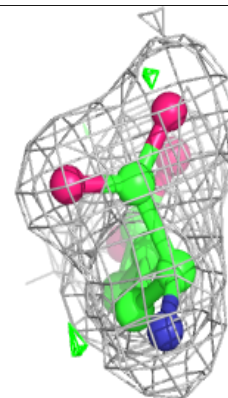
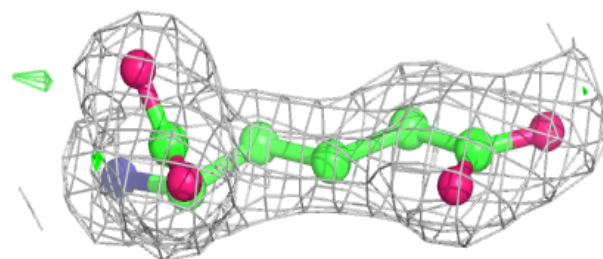
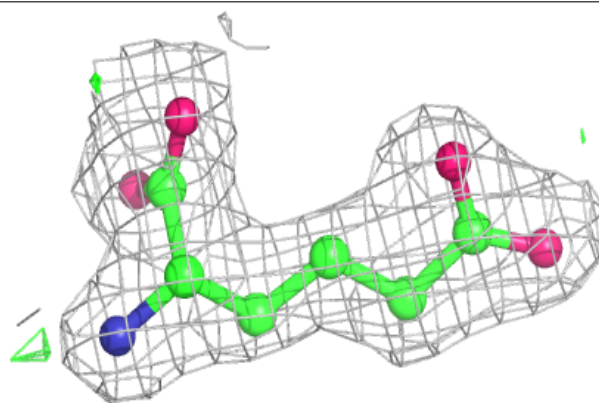


Electron density around UN1 H 601:

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

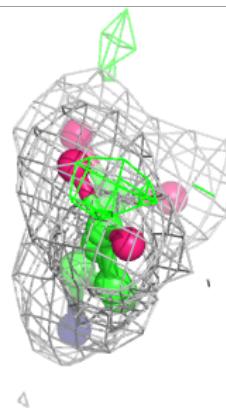
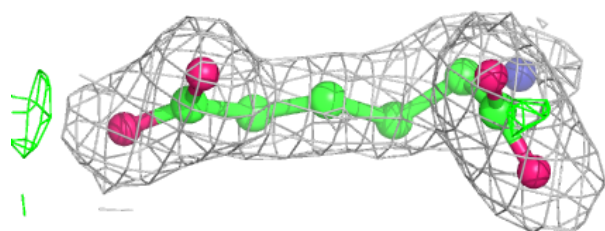
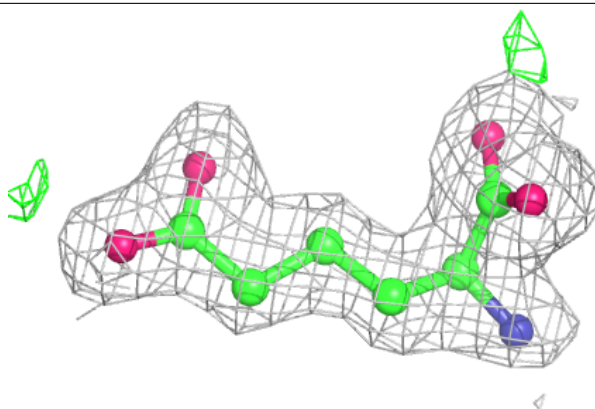
**Electron density around UN1 D 601:**

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and green (positive)

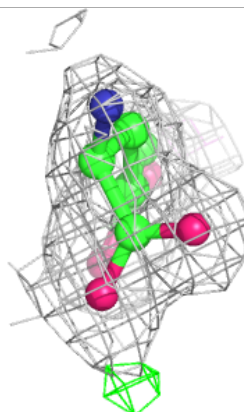
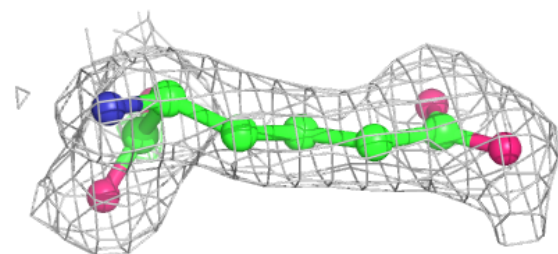
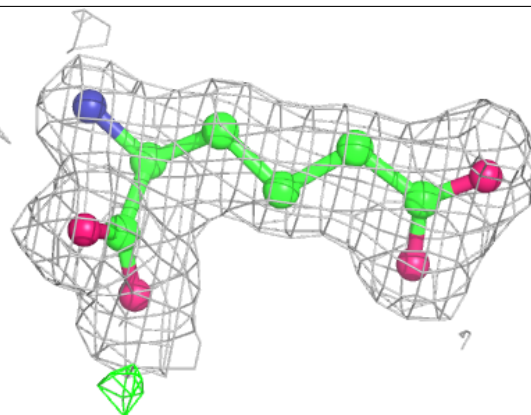


Electron density around UN1 F 601:

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

**Electron density around UN1 B 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.