



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 20, 2026 – 11:50 AM UTC

PDB ID : 6TWM / pdb_00006twm
Title : Product bound structure of the Ectoine utilization protein EutE (DoeB) from *Ruegeria pomeroyi*
Authors : Mais, C.-N.; Altegoer, F.; Bange, G.
Deposited on : 2020-01-13
Resolution : 2.50 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	:	4-5-2 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

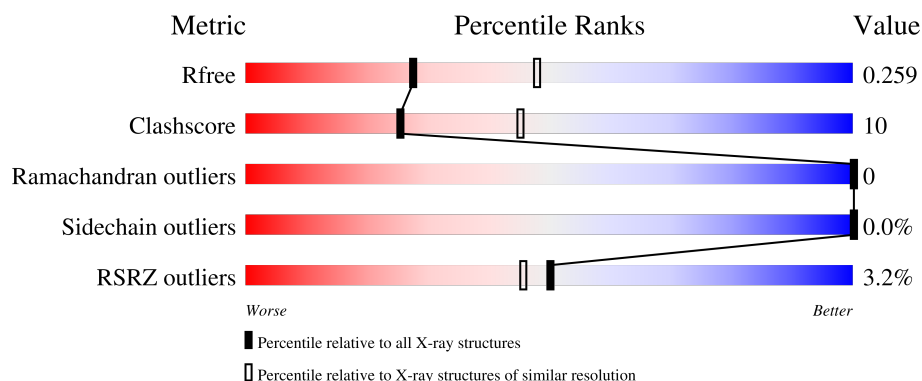
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	337	2% 78% 18% .
1	B	337	4% 78% 19% .
1	C	337	2% 78% 18% .
1	D	337	4% 73% 23% .
1	E	337	0% 79% 18% .

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Mol	Chain	Length	Quality of chain
1	F	337	 % 80% 17% •
1	G	337	 5% 73% 24% •
1	H	337	 3% 80% 17% •
1	I	337	 3% 80% 17% •
1	J	337	 4% 75% 22% •
1	K	337	 2% 81% 16% •
1	L	337	 6% 66% 30% •

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
3	DAB	B	402	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 30865 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 N-acetyl-L-2,4-diaminobutyric acid deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2442	1544	422	464	12			
1	B	326	Total	C	N	O	S	0	0	0
			2442	1544	422	464	12			
1	C	325	Total	C	N	O	S	0	0	0
			2434	1538	421	463	12			
1	D	328	Total	C	N	O	S	0	0	0
			2461	1555	427	467	12			
1	E	327	Total	C	N	O	S	0	0	0
			2451	1549	424	466	12			
1	F	327	Total	C	N	O	S	0	0	0
			2446	1546	423	465	12			
1	G	327	Total	C	N	O	S	0	0	0
			2451	1549	424	466	12			
1	H	327	Total	C	N	O	S	0	0	0
			2451	1549	424	466	12			
1	I	326	Total	C	N	O	S	0	0	0
			2442	1544	422	464	12			
1	J	326	Total	C	N	O	S	0	0	0
			2443	1543	423	465	12			
1	K	327	Total	C	N	O	S	0	0	0
			2451	1549	424	466	12			
1	L	324	Total	C	N	O	S	0	0	0
			2426	1532	421	461	12			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	-	initiating methionine	UNP Q5LUB5
A	-5	GLY	-	expression tag	UNP Q5LUB5
A	-4	HIS	-	expression tag	UNP Q5LUB5
A	-3	HIS	-	expression tag	UNP Q5LUB5
A	-2	HIS	-	expression tag	UNP Q5LUB5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	HIS	-	expression tag	UNP Q5LUB5
A	0	HIS	-	expression tag	UNP Q5LUB5
A	1	HIS	-	expression tag	UNP Q5LUB5
B	-6	MET	-	initiating methionine	UNP Q5LUB5
B	-5	GLY	-	expression tag	UNP Q5LUB5
B	-4	HIS	-	expression tag	UNP Q5LUB5
B	-3	HIS	-	expression tag	UNP Q5LUB5
B	-2	HIS	-	expression tag	UNP Q5LUB5
B	-1	HIS	-	expression tag	UNP Q5LUB5
B	0	HIS	-	expression tag	UNP Q5LUB5
B	1	HIS	-	expression tag	UNP Q5LUB5
C	-6	MET	-	initiating methionine	UNP Q5LUB5
C	-5	GLY	-	expression tag	UNP Q5LUB5
C	-4	HIS	-	expression tag	UNP Q5LUB5
C	-3	HIS	-	expression tag	UNP Q5LUB5
C	-2	HIS	-	expression tag	UNP Q5LUB5
C	-1	HIS	-	expression tag	UNP Q5LUB5
C	0	HIS	-	expression tag	UNP Q5LUB5
C	1	HIS	-	expression tag	UNP Q5LUB5
D	-6	MET	-	initiating methionine	UNP Q5LUB5
D	-5	GLY	-	expression tag	UNP Q5LUB5
D	-4	HIS	-	expression tag	UNP Q5LUB5
D	-3	HIS	-	expression tag	UNP Q5LUB5
D	-2	HIS	-	expression tag	UNP Q5LUB5
D	-1	HIS	-	expression tag	UNP Q5LUB5
D	0	HIS	-	expression tag	UNP Q5LUB5
D	1	HIS	-	expression tag	UNP Q5LUB5
E	-6	MET	-	initiating methionine	UNP Q5LUB5
E	-5	GLY	-	expression tag	UNP Q5LUB5
E	-4	HIS	-	expression tag	UNP Q5LUB5
E	-3	HIS	-	expression tag	UNP Q5LUB5
E	-2	HIS	-	expression tag	UNP Q5LUB5
E	-1	HIS	-	expression tag	UNP Q5LUB5
E	0	HIS	-	expression tag	UNP Q5LUB5
E	1	HIS	-	expression tag	UNP Q5LUB5
F	-6	MET	-	initiating methionine	UNP Q5LUB5
F	-5	GLY	-	expression tag	UNP Q5LUB5
F	-4	HIS	-	expression tag	UNP Q5LUB5
F	-3	HIS	-	expression tag	UNP Q5LUB5
F	-2	HIS	-	expression tag	UNP Q5LUB5
F	-1	HIS	-	expression tag	UNP Q5LUB5
F	0	HIS	-	expression tag	UNP Q5LUB5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	HIS	-	expression tag	UNP Q5LUB5
G	-6	MET	-	initiating methionine	UNP Q5LUB5
G	-5	GLY	-	expression tag	UNP Q5LUB5
G	-4	HIS	-	expression tag	UNP Q5LUB5
G	-3	HIS	-	expression tag	UNP Q5LUB5
G	-2	HIS	-	expression tag	UNP Q5LUB5
G	-1	HIS	-	expression tag	UNP Q5LUB5
G	0	HIS	-	expression tag	UNP Q5LUB5
G	1	HIS	-	expression tag	UNP Q5LUB5
H	-6	MET	-	initiating methionine	UNP Q5LUB5
H	-5	GLY	-	expression tag	UNP Q5LUB5
H	-4	HIS	-	expression tag	UNP Q5LUB5
H	-3	HIS	-	expression tag	UNP Q5LUB5
H	-2	HIS	-	expression tag	UNP Q5LUB5
H	-1	HIS	-	expression tag	UNP Q5LUB5
H	0	HIS	-	expression tag	UNP Q5LUB5
H	1	HIS	-	expression tag	UNP Q5LUB5
I	-6	MET	-	initiating methionine	UNP Q5LUB5
I	-5	GLY	-	expression tag	UNP Q5LUB5
I	-4	HIS	-	expression tag	UNP Q5LUB5
I	-3	HIS	-	expression tag	UNP Q5LUB5
I	-2	HIS	-	expression tag	UNP Q5LUB5
I	-1	HIS	-	expression tag	UNP Q5LUB5
I	0	HIS	-	expression tag	UNP Q5LUB5
I	1	HIS	-	expression tag	UNP Q5LUB5
J	-6	MET	-	initiating methionine	UNP Q5LUB5
J	-5	GLY	-	expression tag	UNP Q5LUB5
J	-4	HIS	-	expression tag	UNP Q5LUB5
J	-3	HIS	-	expression tag	UNP Q5LUB5
J	-2	HIS	-	expression tag	UNP Q5LUB5
J	-1	HIS	-	expression tag	UNP Q5LUB5
J	0	HIS	-	expression tag	UNP Q5LUB5
J	1	HIS	-	expression tag	UNP Q5LUB5
K	-6	MET	-	initiating methionine	UNP Q5LUB5
K	-5	GLY	-	expression tag	UNP Q5LUB5
K	-4	HIS	-	expression tag	UNP Q5LUB5
K	-3	HIS	-	expression tag	UNP Q5LUB5
K	-2	HIS	-	expression tag	UNP Q5LUB5
K	-1	HIS	-	expression tag	UNP Q5LUB5
K	0	HIS	-	expression tag	UNP Q5LUB5
K	1	HIS	-	expression tag	UNP Q5LUB5
L	-6	MET	-	initiating methionine	UNP Q5LUB5

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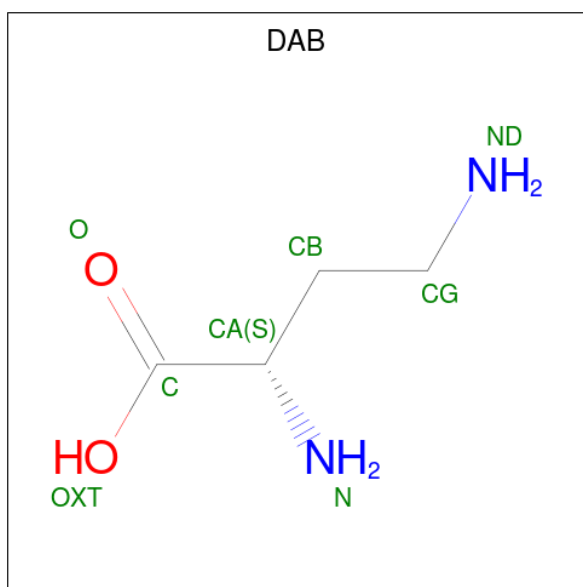
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Chain	Residue	Modelled	Actual	Comment	Reference
L	-5	GLY	-	expression tag	UNP Q5LUB5
L	-4	HIS	-	expression tag	UNP Q5LUB5
L	-3	HIS	-	expression tag	UNP Q5LUB5
L	-2	HIS	-	expression tag	UNP Q5LUB5
L	-1	HIS	-	expression tag	UNP Q5LUB5
L	0	HIS	-	expression tag	UNP Q5LUB5
L	1	HIS	-	expression tag	UNP Q5LUB5

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

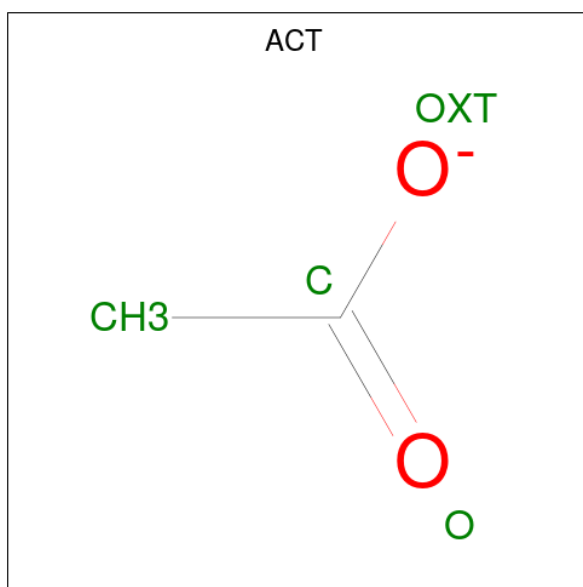
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	D	1	Total 1	Zn 1	0	0
2	E	1	Total 1	Zn 1	0	0
2	G	1	Total 1	Zn 1	0	0
2	H	1	Total 1	Zn 1	0	0
2	I	1	Total 1	Zn 1	0	0
2	J	1	Total 1	Zn 1	0	0
2	L	1	Total 1	Zn 1	0	0

- Molecule 3 is 2,4-DIAMINOBTYRIC ACID (CCD ID: DAB) (formula: C₄H₁₀N₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	2	2		
3	B	1	Total	C	N	O	0	0
			8	4	2	2		
3	D	1	Total	C	N	O	0	0
			8	4	2	2		
3	E	1	Total	C	N	O	0	0
			8	4	2	2		
3	F	1	Total	C	N	O	0	0
			8	4	2	2		
3	J	1	Total	C	N	O	0	0
			8	4	2	2		
3	K	1	Total	C	N	O	0	0
			8	4	2	2		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	144	Total	O	0	0
			144	144		
5	B	130	Total	O	0	0
			130	130		
5	C	130	Total	O	0	0
			130	130		
5	D	111	Total	O	0	0
			111	111		
5	E	117	Total	O	0	0
			117	117		
5	F	131	Total	O	0	0
			131	131		
5	G	121	Total	O	0	0
			121	121		
5	H	112	Total	O	0	0
			112	112		
5	I	108	Total	O	0	0
			108	108		

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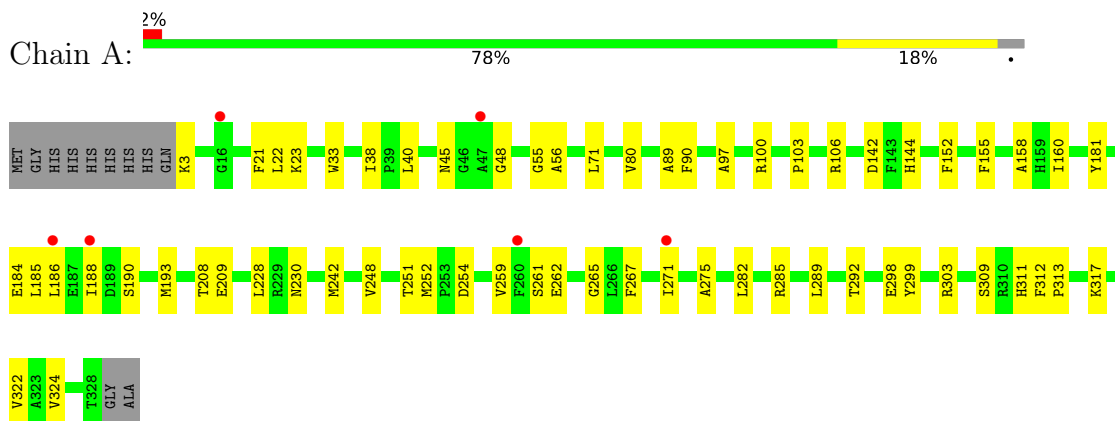
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	118	Total 118	O 118	0	0
5	K	118	Total 118	O 118	0	0
5	L	107	Total 107	O 107	0	0

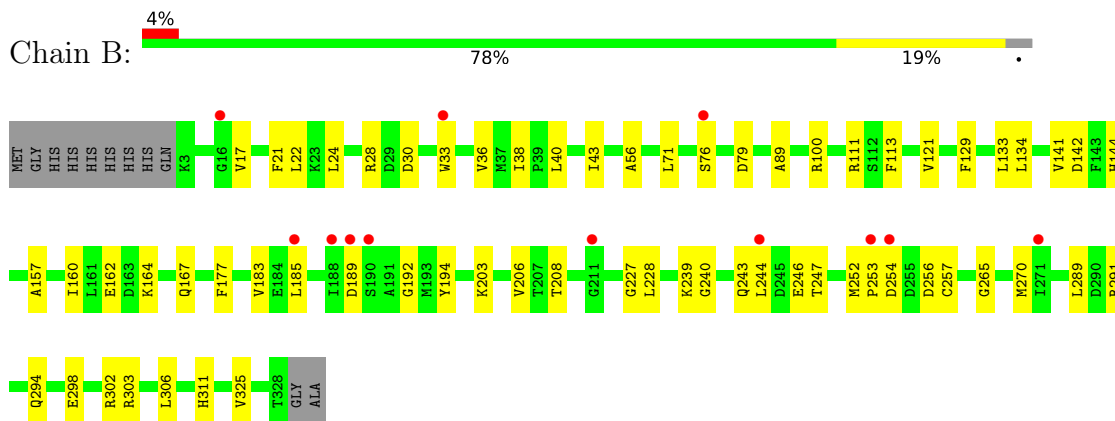
3 Residue-property plots [i](#)

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.

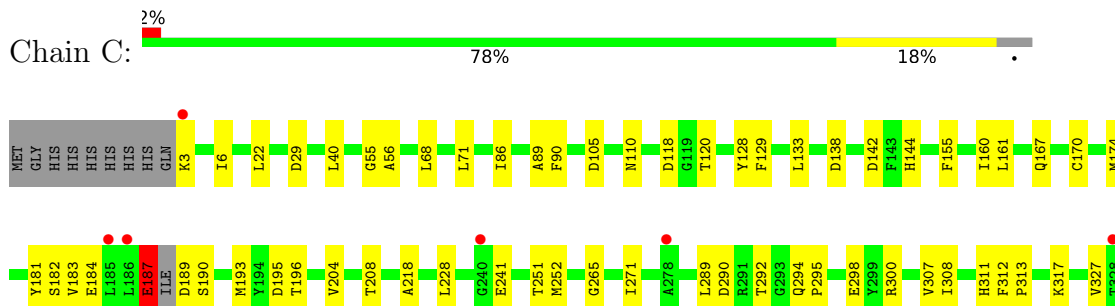
- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase



- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

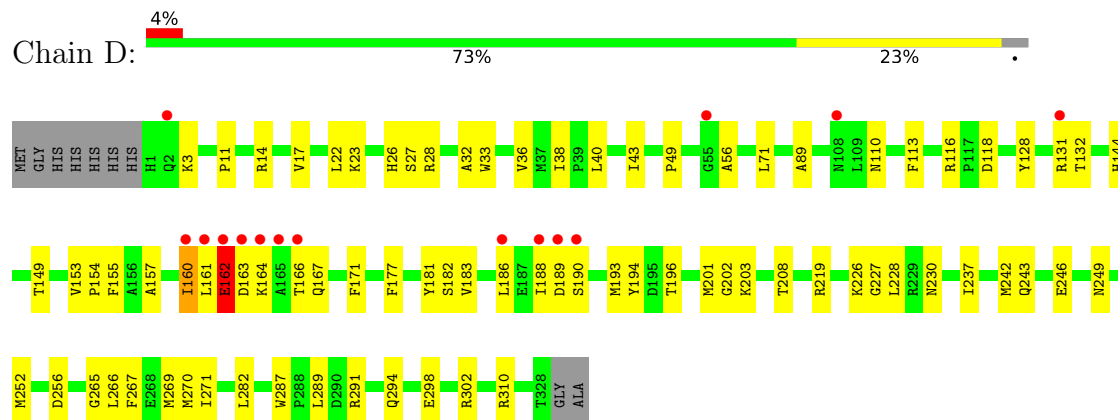


- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

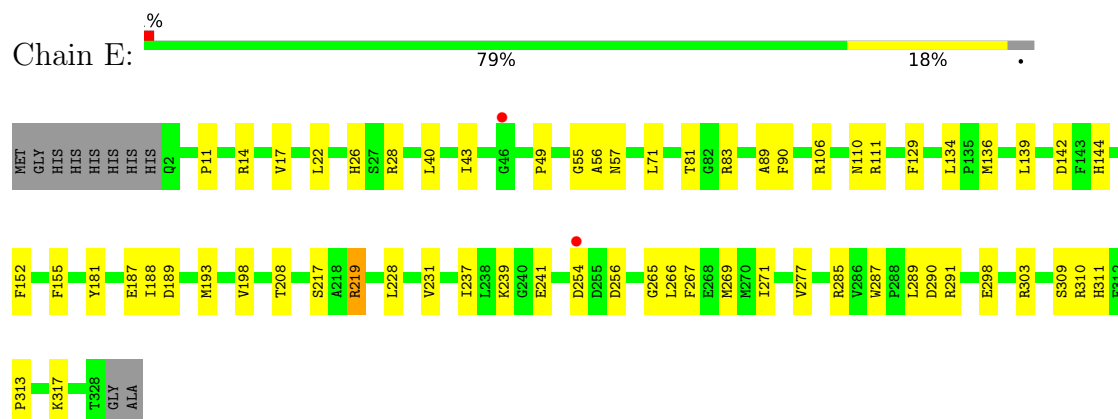


GLY
ALA

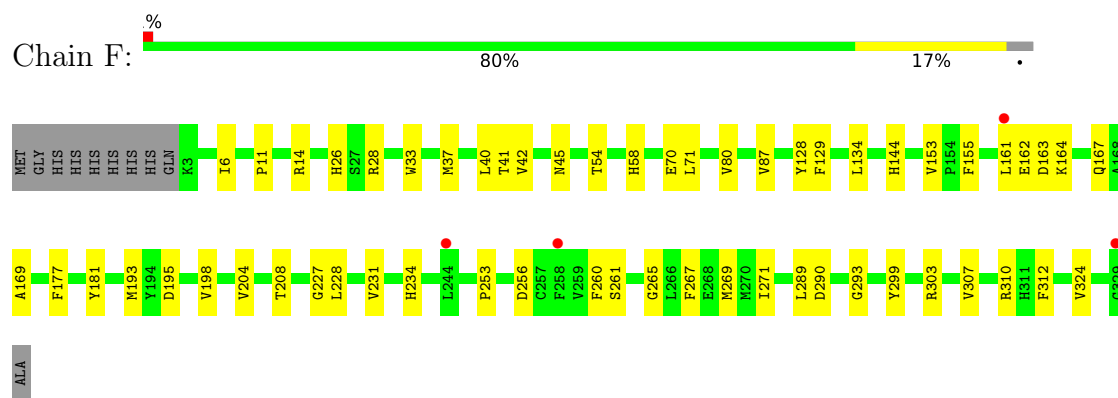
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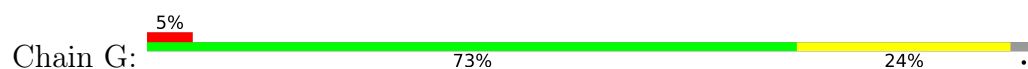


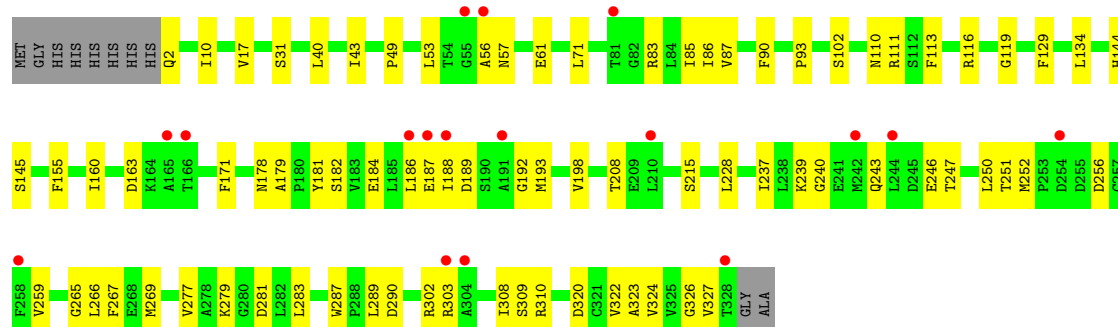
- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase



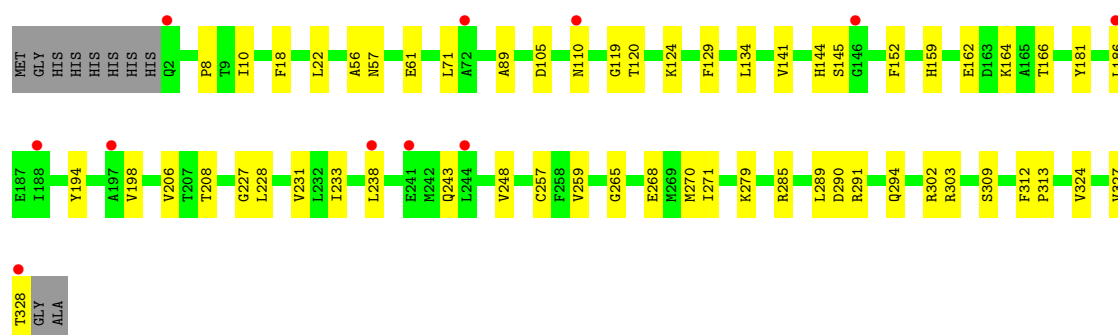
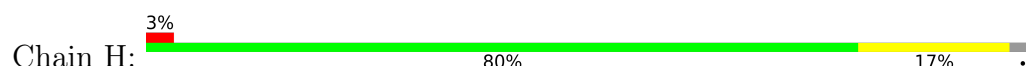
ALA

- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

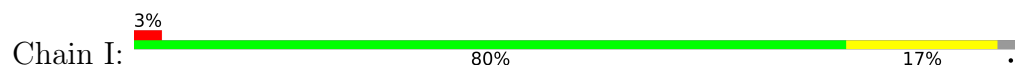




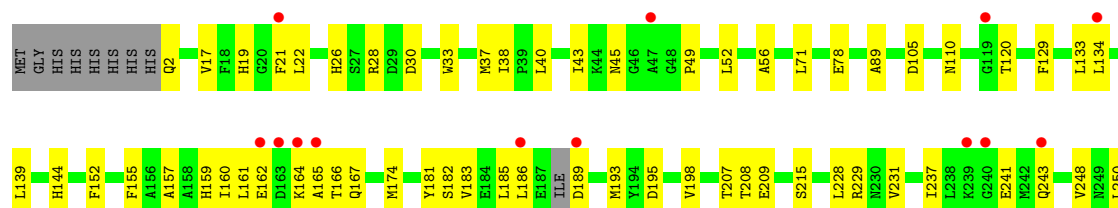
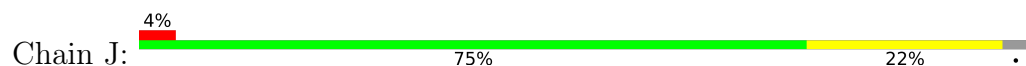
- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

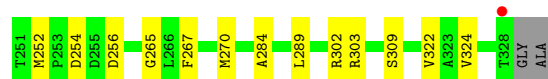


- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

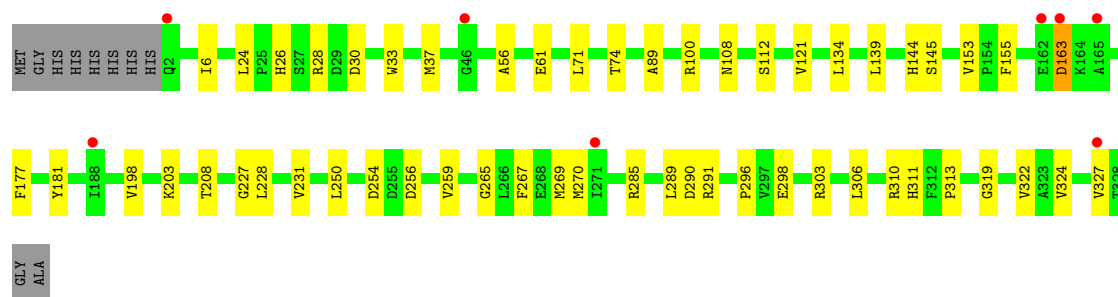
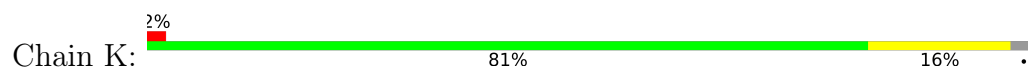


- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase

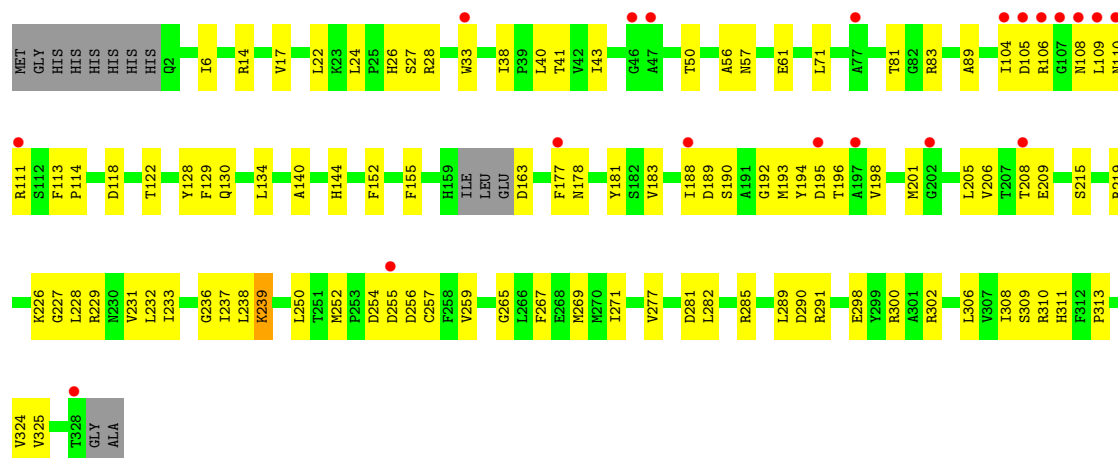




- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase



- Molecule 1: N-acetyl-L-2,4-diaminobutyric acid deacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.67Å 145.95Å 164.12Å 90.00° 92.30° 90.00°	Depositor
Resolution (Å)	46.64 – 2.50 46.64 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (46.64-2.50) 98.8 (46.64-2.50)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.209 , 0.262 0.207 , 0.259	Depositor DCC
R_{free} test set	1568 reflections (0.94%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	30865	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.38 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.8815e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: DAB, ZN, ACT

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	0.14	0/2495	0.38	0/3390
1	B	0.15	0/2495	0.41	0/3390
1	C	0.18	0/2487	0.44	1/3379 (0.0%)
1	D	0.26	2/2515 (0.1%)	0.50	5/3417 (0.1%)
1	E	0.17	0/2504	0.40	1/3402 (0.0%)
1	F	0.15	0/2499	0.39	0/3395
1	G	0.16	0/2504	0.44	0/3402
1	H	0.15	0/2504	0.41	0/3402
1	I	0.22	1/2495 (0.0%)	0.46	0/3390
1	J	0.16	0/2495	0.40	0/3388
1	K	0.18	0/2504	0.43	1/3402 (0.0%)
1	L	0.16	0/2478	0.43	0/3365
All	All	0.18	3/29975 (0.0%)	0.42	8/40722 (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	C	0	1
1	D	0	3
1	G	0	1
1	J	0	1
1	K	0	1
1	L	0	1
All	All	0	9

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	131	ARG	NE-CZ	7.71	1.41	1.33
1	I	106	ARG	CD-NE	6.20	1.54	1.46
1	D	131	ARG	CZ-NH1	-5.56	1.25	1.32

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	131	ARG	NE-CZ-NH2	7.01	125.51	119.20
1	D	162	GLU	CA-C-N	6.16	132.22	122.61
1	D	162	GLU	C-N-CA	6.16	132.22	122.61
1	K	163	ASP	CA-CB-CG	-5.63	106.97	112.60
1	D	162	GLU	N-CA-C	5.48	119.12	109.80

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	ILE	Peptide
1	C	187	GLU	Peptide
1	D	160	ILE	Peptide
1	D	161	LEU	Peptide
1	D	162	GLU	Peptide

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	2442	0	2423	45	0
1	B	2442	0	2423	45	3
1	C	2434	0	2412	40	1
1	D	2461	0	2441	57	0
1	E	2451	0	2431	42	0
1	F	2446	0	2426	42	0
1	G	2451	0	2431	63	0
1	H	2451	0	2431	37	0
1	I	2442	0	2423	42	1
1	J	2443	0	2419	52	0
1	K	2451	0	2431	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	2426	0	2402	81	4
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	L	1	0	0	0	0
3	A	8	0	8	1	0
3	B	8	0	9	2	3
3	D	8	0	9	0	0
3	E	8	0	9	1	2
3	F	8	0	9	0	0
3	J	8	0	9	1	0
3	K	8	0	9	2	0
4	C	4	0	3	0	0
4	G	4	0	3	0	0
4	H	4	0	3	0	0
5	A	144	0	0	6	0
5	B	130	0	0	4	0
5	C	130	0	0	7	0
5	D	111	0	0	4	0
5	E	117	0	0	4	0
5	F	131	0	0	2	0
5	G	121	0	0	10	0
5	H	112	0	0	4	0
5	I	108	0	0	5	0
5	J	118	0	0	7	0
5	K	118	0	0	0	0
5	L	107	0	0	9	0
All	All	30865	0	29164	564	7

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 10.

The worst 5 of 564 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:22:LEU:HG	1:B:40:LEU:HD22	1.28	1.06
1:I:22:LEU:HG	1:I:40:LEU:HD22	1.34	1.05
1:I:103:PRO:HA	1:I:106:ARG:HH21	1.23	1.03
1:D:22:LEU:HG	1:D:40:LEU:HD22	1.40	1.00
1:G:303:ARG:NH1	1:G:303:ARG:HB3	1.80	0.97

The worst 5 of 7 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)
1:L:33:TRP:CE2	3:B:402:DAB:ND[1_545]	1.44	0.76
1:B:33:TRP:CZ2	3:E:402:DAB:CG[2_555]	1.49	0.71
1:L:33:TRP:CZ2	3:B:402:DAB:ND[1_545]	1.66	0.54
1:B:100:ARG:NH1	1:L:27:SER:OG[1_565]	2.08	0.12
1:L:33:TRP:NE1	3:B:402:DAB:ND[1_545]	2.10	0.10

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	324/337 (96%)	318 (98%)	6 (2%)	0	100	100
1	B	324/337 (96%)	318 (98%)	6 (2%)	0	100	100
1	C	323/337 (96%)	315 (98%)	8 (2%)	0	100	100
1	D	326/337 (97%)	314 (96%)	12 (4%)	0	100	100
1	E	325/337 (96%)	322 (99%)	3 (1%)	0	100	100
1	F	325/337 (96%)	319 (98%)	6 (2%)	0	100	100
1	G	325/337 (96%)	316 (97%)	9 (3%)	0	100	100
1	H	325/337 (96%)	319 (98%)	6 (2%)	0	100	100
1	I	324/337 (96%)	320 (99%)	4 (1%)	0	100	100
1	J	322/337 (96%)	318 (99%)	4 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	325/337 (96%)	320 (98%)	5 (2%)	0	100	100
1	L	320/337 (95%)	308 (96%)	12 (4%)	0	100	100
All	All	3888/4044 (96%)	3807 (98%)	81 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	257/265 (97%)	257 (100%)	0	100	100
1	B	257/265 (97%)	257 (100%)	0	100	100
1	C	256/265 (97%)	256 (100%)	0	100	100
1	D	259/265 (98%)	259 (100%)	0	100	100
1	E	258/265 (97%)	258 (100%)	0	100	100
1	F	257/265 (97%)	257 (100%)	0	100	100
1	G	258/265 (97%)	258 (100%)	0	100	100
1	H	258/265 (97%)	257 (100%)	1 (0%)	84	93
1	I	257/265 (97%)	257 (100%)	0	100	100
1	J	257/265 (97%)	257 (100%)	0	100	100
1	K	258/265 (97%)	258 (100%)	0	100	100
1	L	255/265 (96%)	255 (100%)	0	100	100
All	All	3087/3180 (97%)	3086 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	328	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	K	69	HIS
1	L	221	ASN
1	I	69	HIS
1	I	130	GLN
1	I	159	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 20 ligands modelled in this entry, 10 are monoatomic - leaving 10 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	$\# Z > 2$
3	DAB	D	402	2	6,7,7	0.88	0	4,8,8	0.74	0
3	DAB	B	402	2	6,7,7	0.87	0	4,8,8	1.14	1 (25%)
3	DAB	A	402	2,1	6,7,7	0.88	0	4,8,8	0.71	0
3	DAB	E	402	2	6,7,7	0.85	0	4,8,8	0.76	0
4	ACT	H	402	2	3,3,3	1.37	0	3,3,3	1.47	0
3	DAB	F	401	-	6,7,7	0.86	0	4,8,8	0.69	0
4	ACT	C	402	2	3,3,3	1.29	0	3,3,3	1.53	0
3	DAB	K	401	-	6,7,7	0.85	0	4,8,8	0.60	0
3	DAB	J	402	2	6,7,7	0.82	0	4,8,8	0.55	0
4	ACT	G	402	-	3,3,3	1.40	1 (33%)	3,3,3	1.35	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
3	DAB	D	402	2	-	3/7/7/7	-
3	DAB	B	402	2	-	3/7/7/7	-
3	DAB	A	402	2,1	-	1/7/7/7	-
3	DAB	E	402	2	-	0/7/7/7	-
3	DAB	F	401	-	-	0/7/7/7	-
3	DAB	K	401	-	-	5/7/7/7	-
3	DAB	J	402	2	-	5/7/7/7	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	402	ACT	CH3-C	2.02	1.57	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	DAB	CB-CA-C	-2.06	105.00	110.45

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	402	DAB	CA-CB-CG-ND
3	J	402	DAB	O-C-CA-N
3	J	402	DAB	CA-CB-CG-ND
3	K	401	DAB	O-C-CA-N
3	J	402	DAB	OXT-C-CA-N

There are no ring outliers.

5 monomers are involved in 12 short contacts:

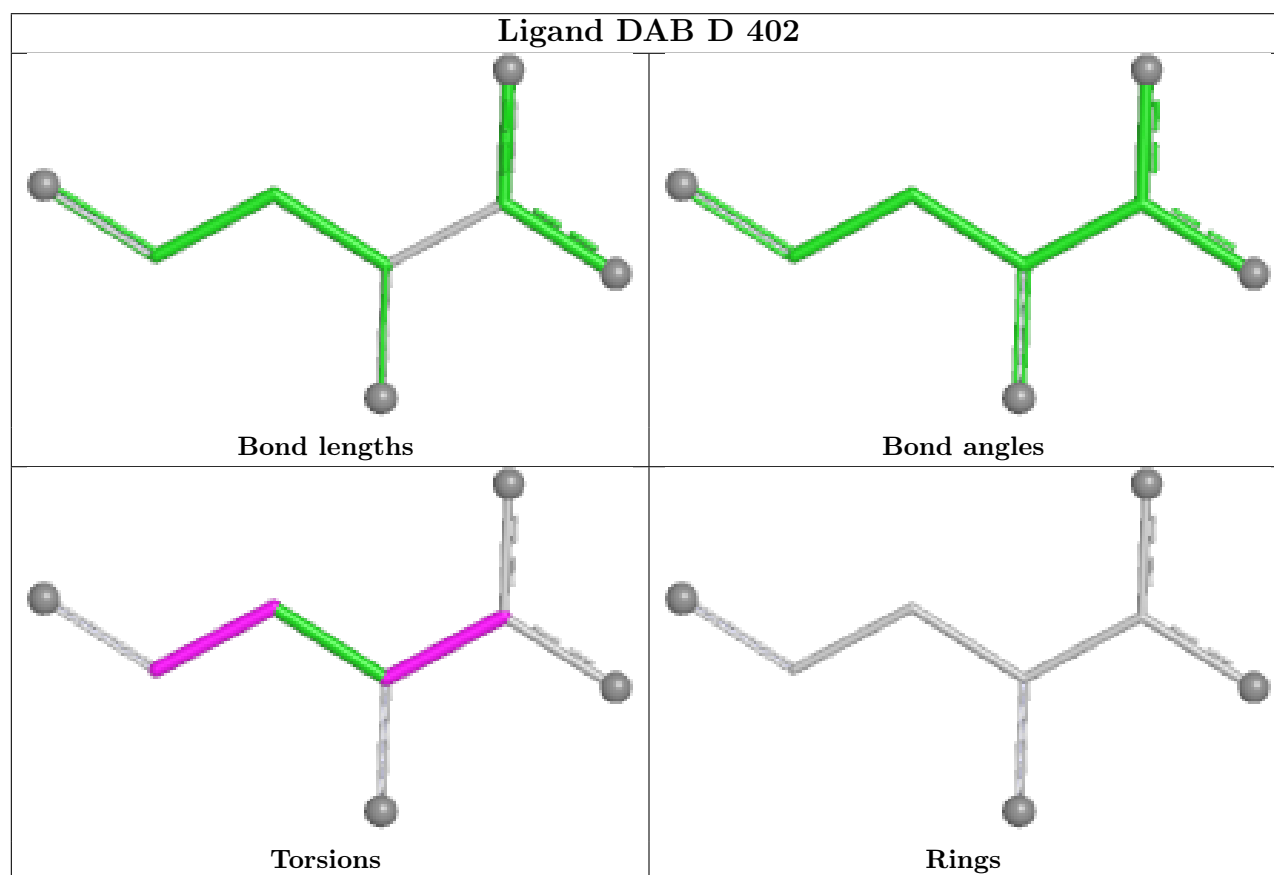
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	DAB	2	3
3	A	402	DAB	1	0
3	E	402	DAB	1	2
3	K	401	DAB	2	0

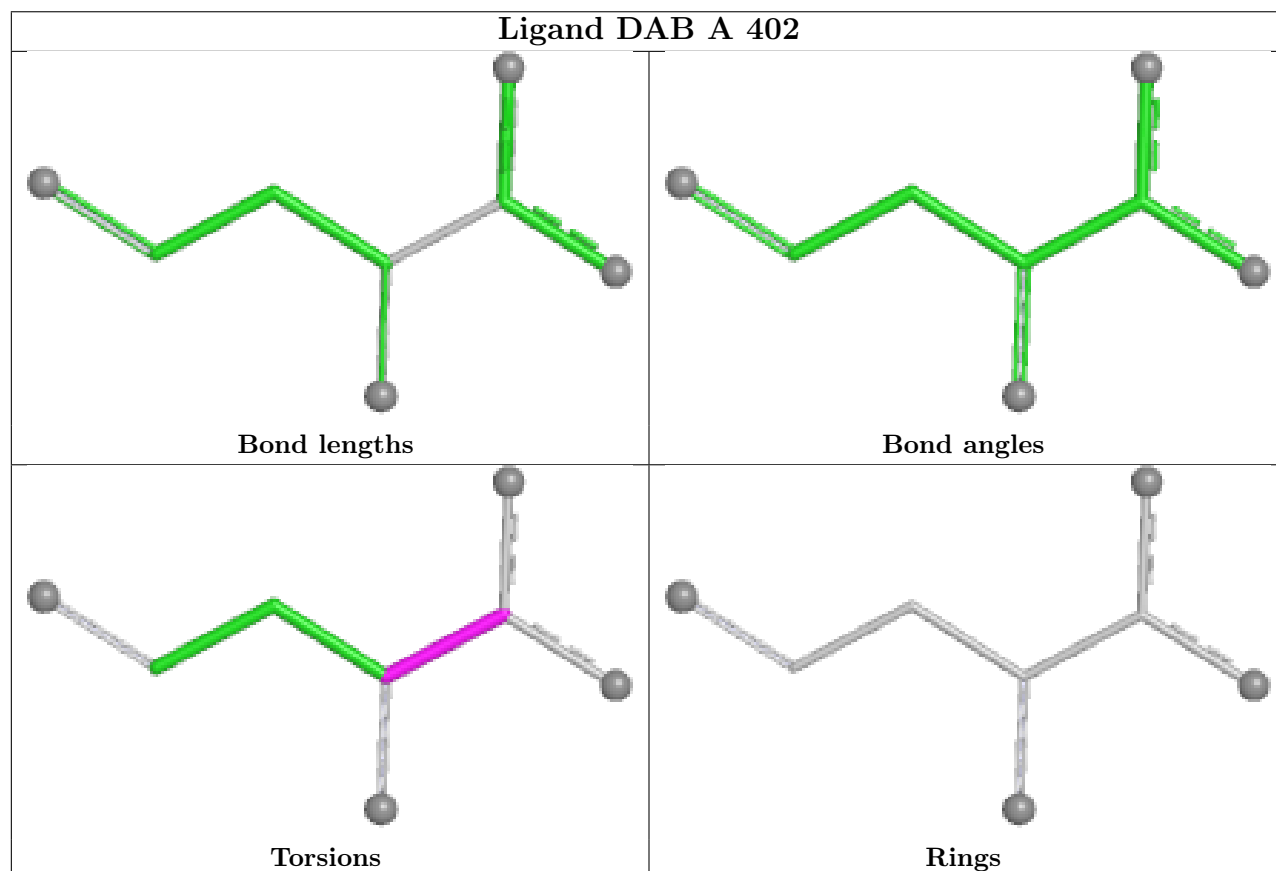
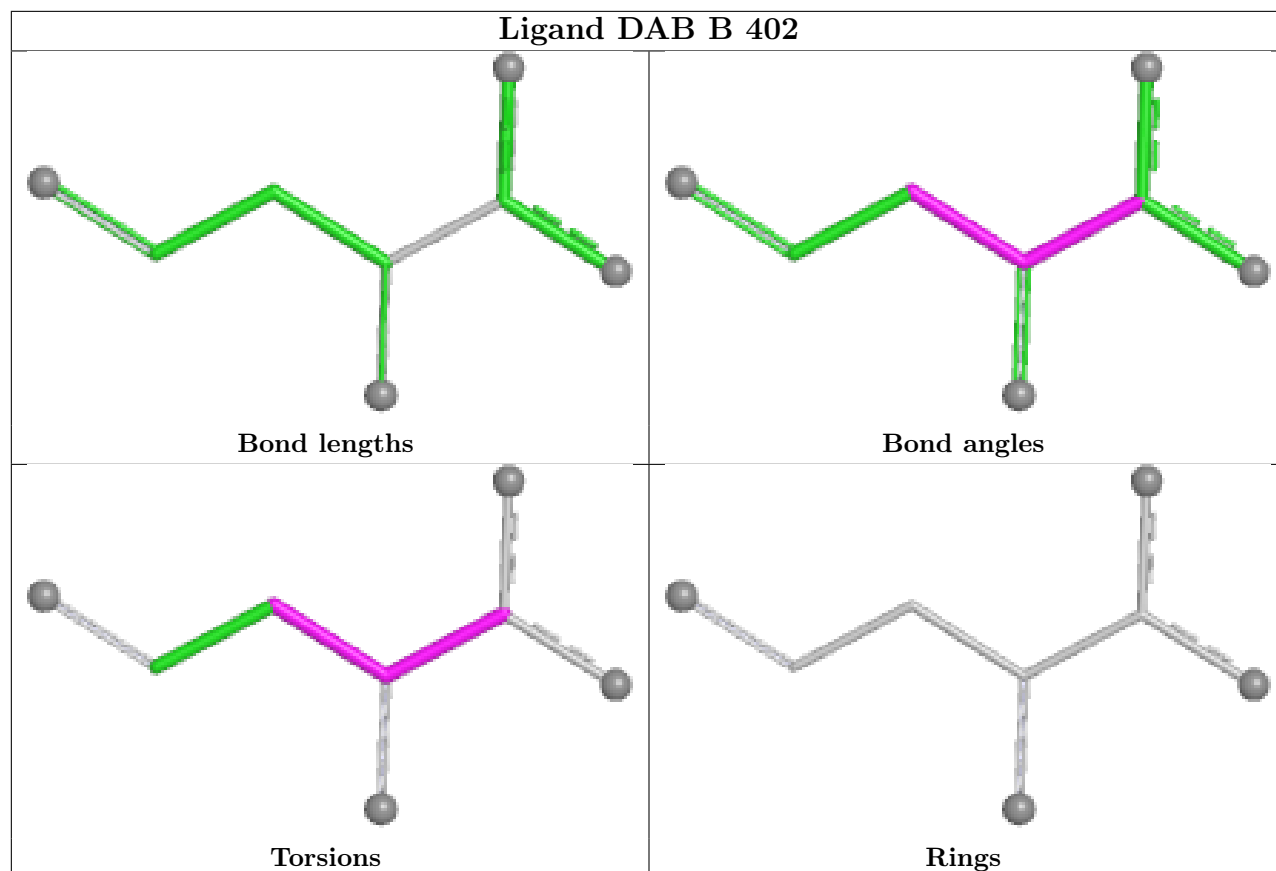
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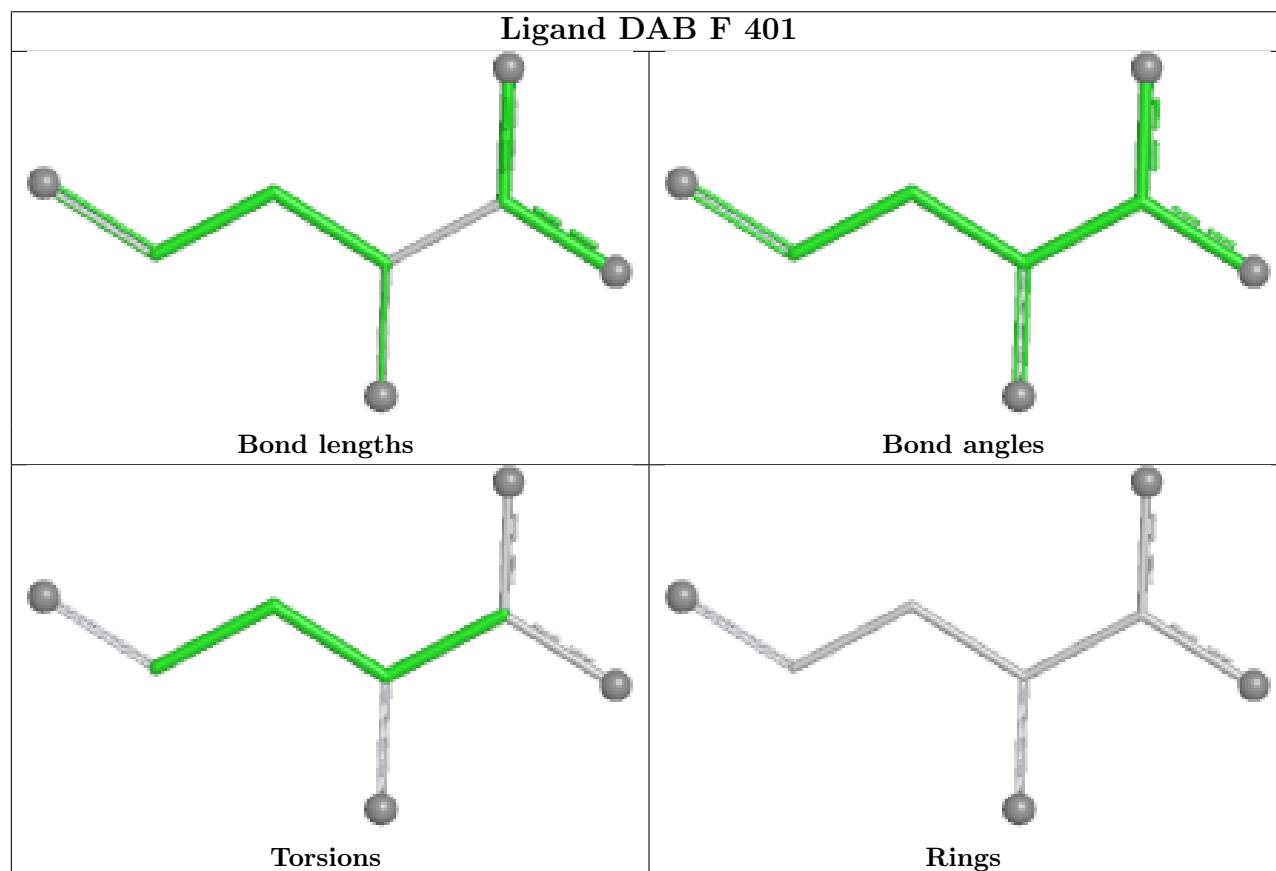
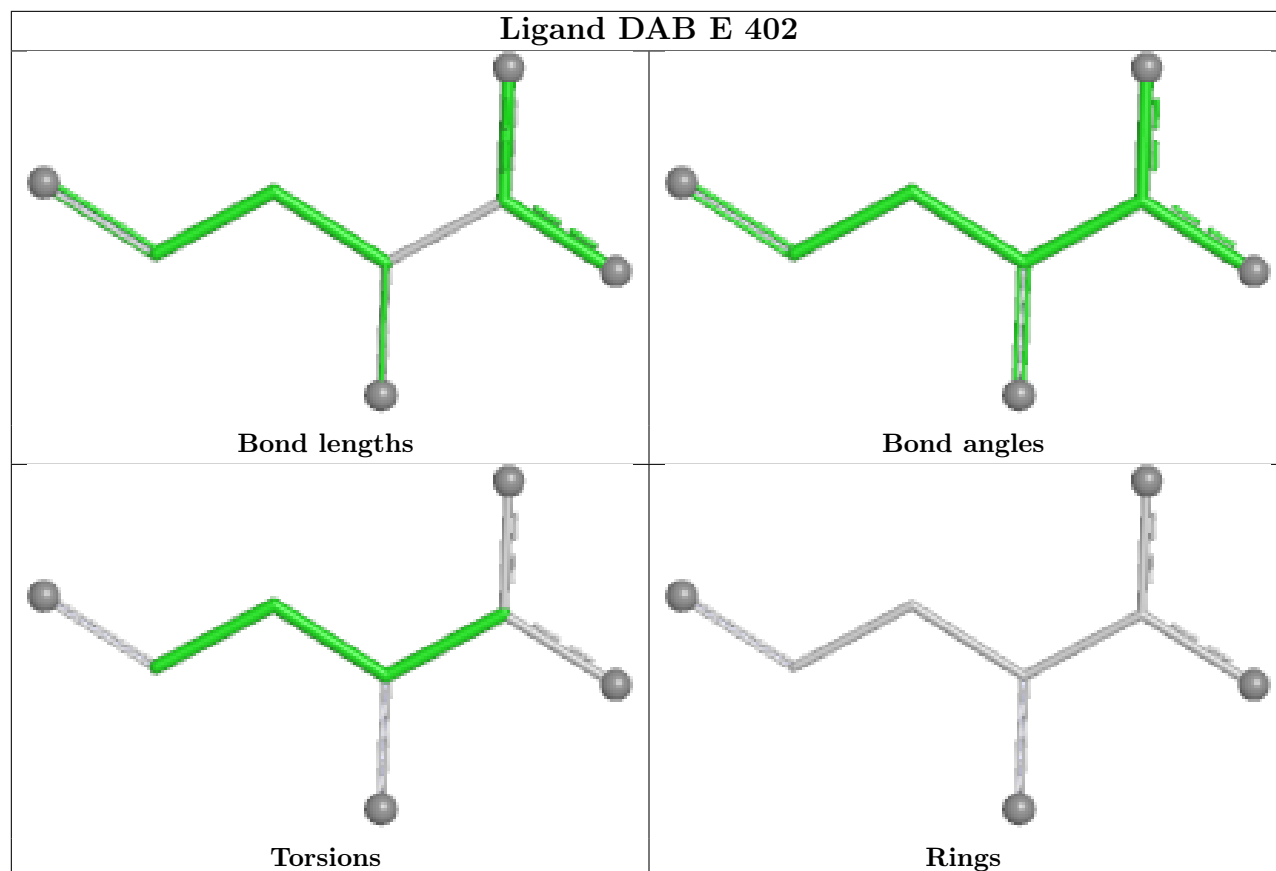
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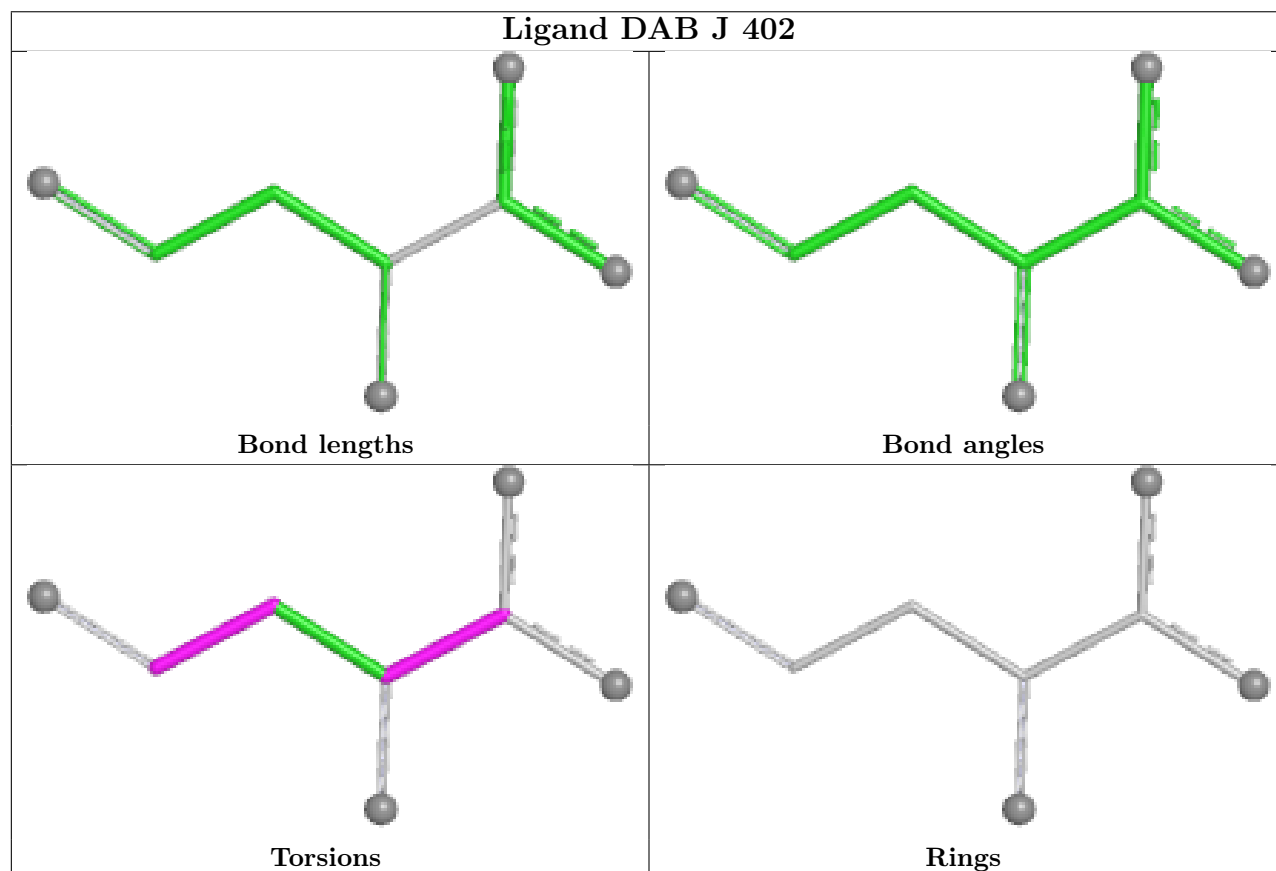
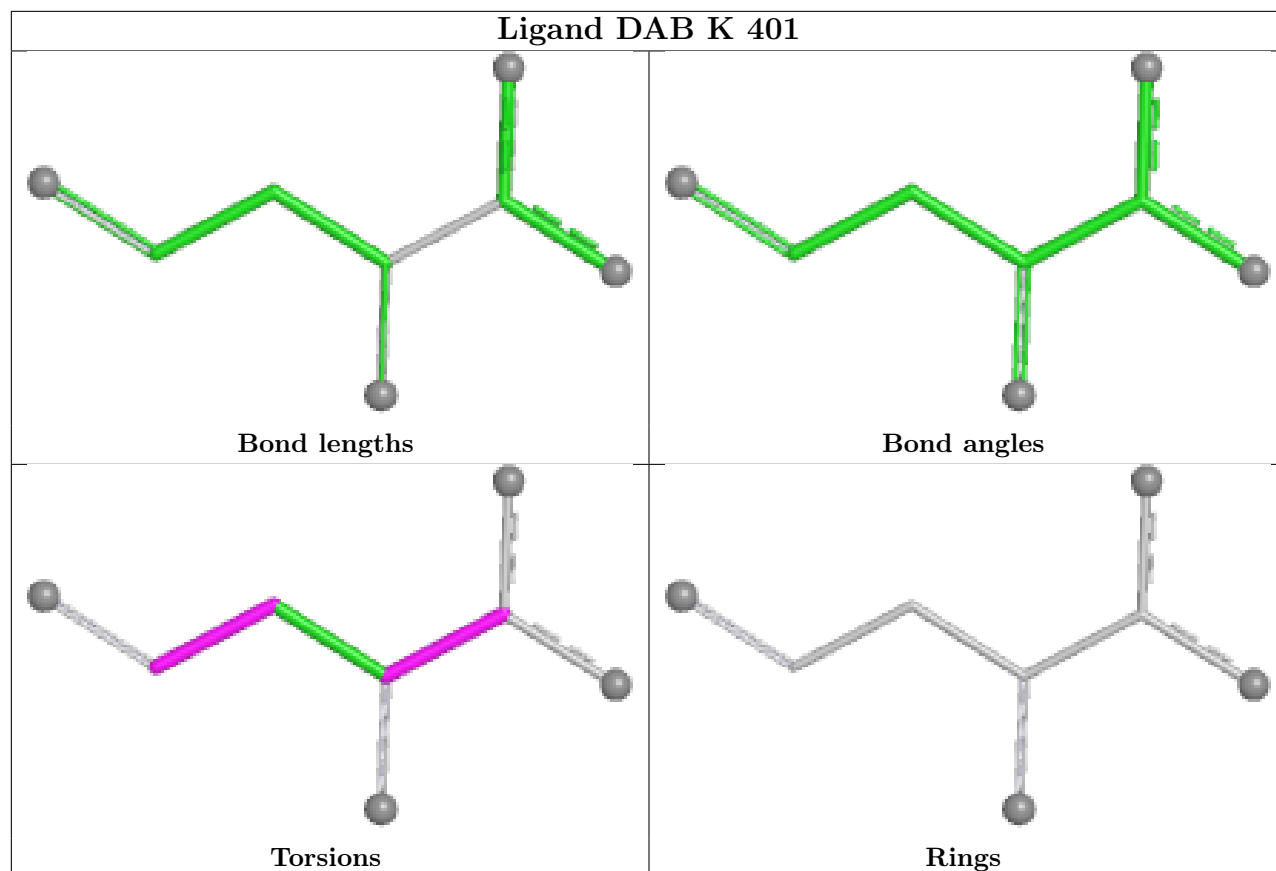
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	402	DAB	1	0

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	326/337 (96%)	0.18	6 (1%) 67 64	28, 41, 72, 125	0
1	B	326/337 (96%)	0.32	12 (3%) 45 40	27, 44, 74, 139	0
1	C	325/337 (96%)	0.27	6 (1%) 67 64	24, 44, 73, 115	0
1	D	328/337 (97%)	0.29	15 (4%) 37 33	25, 43, 76, 132	0
1	E	327/337 (97%)	0.19	2 (0%) 85 83	28, 43, 68, 95	0
1	F	327/337 (97%)	0.22	4 (1%) 76 73	26, 44, 66, 92	0
1	G	327/337 (97%)	0.50	17 (5%) 33 29	30, 48, 83, 117	0
1	H	327/337 (97%)	0.38	11 (3%) 48 43	31, 48, 84, 125	0
1	I	326/337 (96%)	0.44	9 (2%) 55 50	29, 51, 88, 139	0
1	J	326/337 (96%)	0.36	14 (4%) 40 35	29, 47, 83, 138	0
1	K	327/337 (97%)	0.42	8 (2%) 59 55	31, 48, 76, 130	0
1	L	324/337 (96%)	0.75	20 (6%) 26 23	34, 60, 92, 130	0
All	All	3916/4044 (96%)	0.36	124 (3%) 50 46	24, 46, 80, 139	0

The worst 5 of 124 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	163	ASP	8.2
1	L	107	GLY	6.7
1	L	108	ASN	6.0
1	J	119	GLY	5.4
1	C	240	GLY	4.9

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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

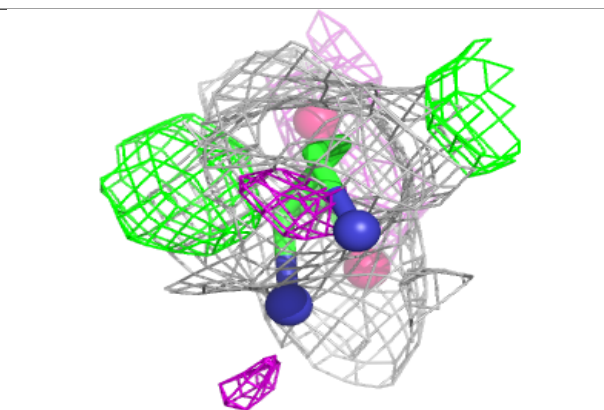
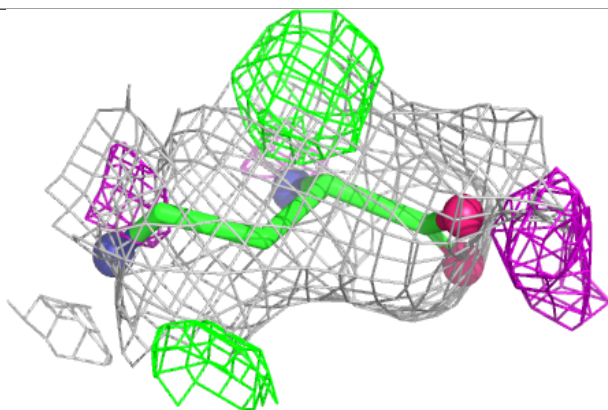
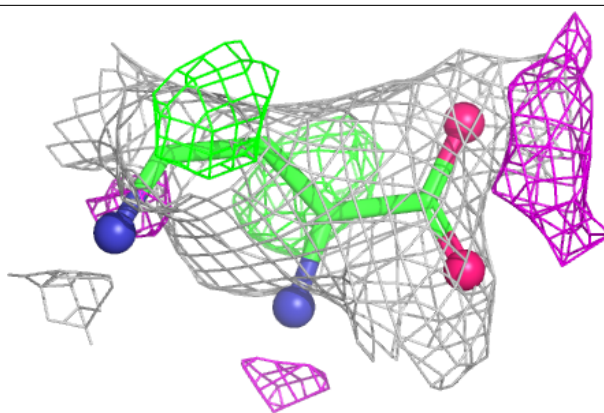
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	DAB	E	402	8/8	0.55	0.24	59,77,89,95	0
4	ACT	G	402	4/4	0.55	0.29	82,83,83,83	0
4	ACT	H	402	4/4	0.69	0.16	75,75,76,76	0
3	DAB	D	402	8/8	0.72	0.24	84,85,86,86	0
2	ZN	G	401	1/1	0.75	0.29	83,83,83,83	0
3	DAB	B	402	8/8	0.75	0.18	59,76,78,81	0
3	DAB	K	401	8/8	0.77	0.22	90,91,93,93	0
3	DAB	F	401	8/8	0.79	0.16	61,62,63,64	0
2	ZN	D	401	1/1	0.79	0.18	134,134,134,134	0
4	ACT	C	402	4/4	0.80	0.28	110,110,111,112	0
2	ZN	L	401	1/1	0.80	0.12	133,133,133,133	0
2	ZN	H	401	1/1	0.80	0.17	132,132,132,132	0
3	DAB	A	402	8/8	0.82	0.21	100,101,105,105	0
3	DAB	J	402	8/8	0.82	0.20	85,87,89,90	0
2	ZN	C	401	1/1	0.86	0.17	114,114,114,114	0
2	ZN	E	401	1/1	0.88	0.24	101,101,101,101	0
2	ZN	J	401	1/1	0.93	0.17	72,72,72,72	0
2	ZN	B	401	1/1	0.93	0.16	84,84,84,84	0
2	ZN	A	401	1/1	0.96	0.12	101,101,101,101	0
2	ZN	I	401	1/1	0.96	0.10	95,95,95,95	0

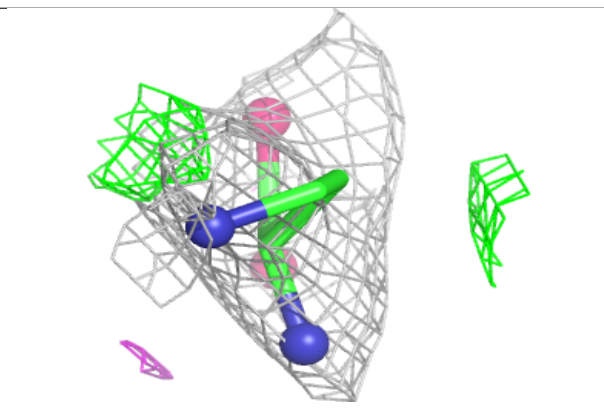
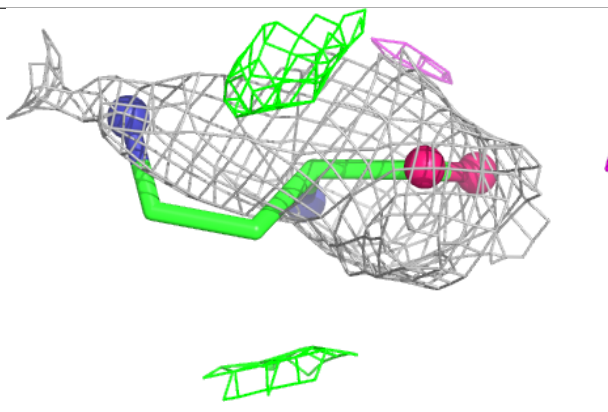
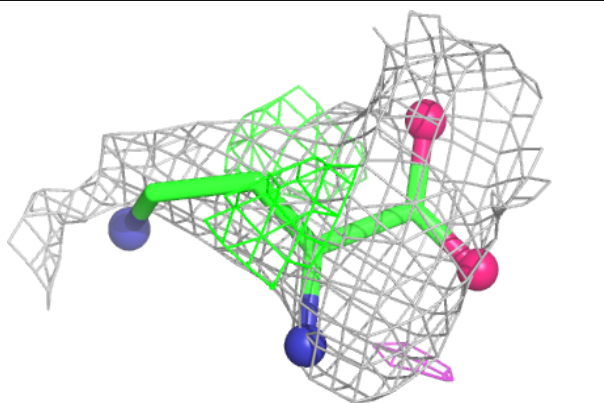
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DAB E 402:

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

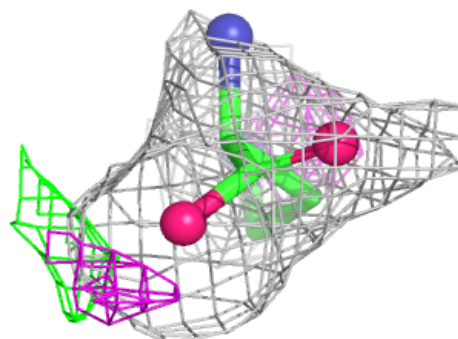
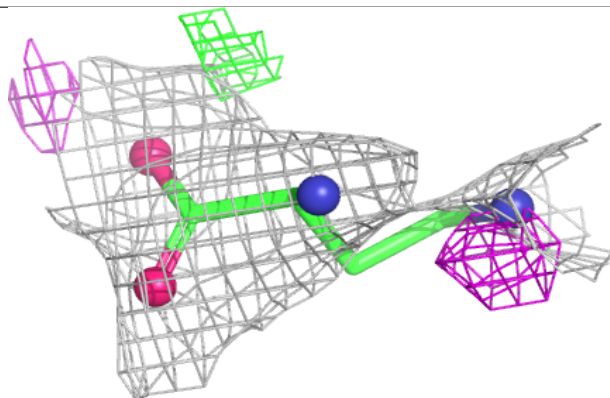
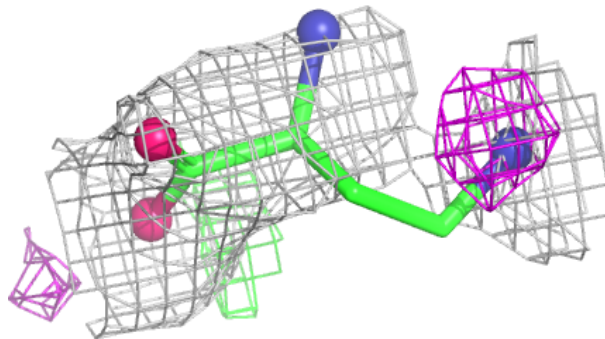
**Electron density around DAB D 402:**

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

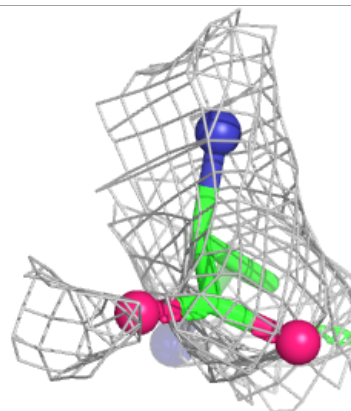
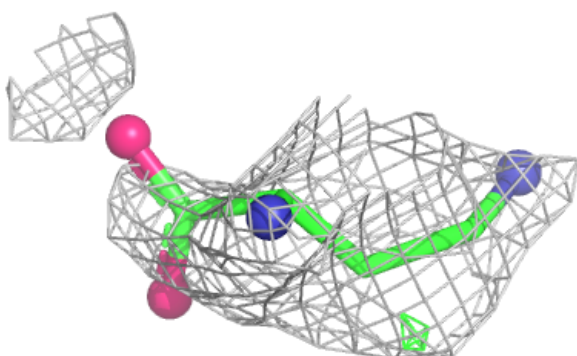
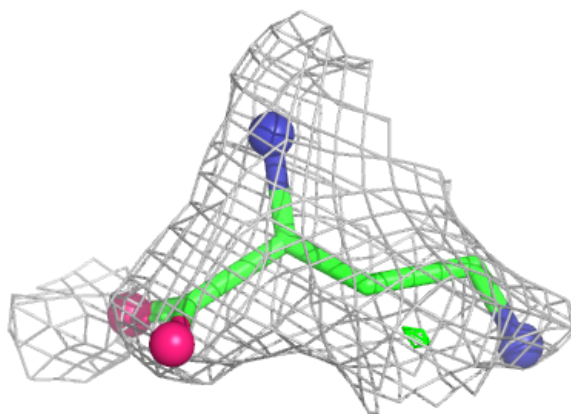


Electron density around DAB B 402:

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

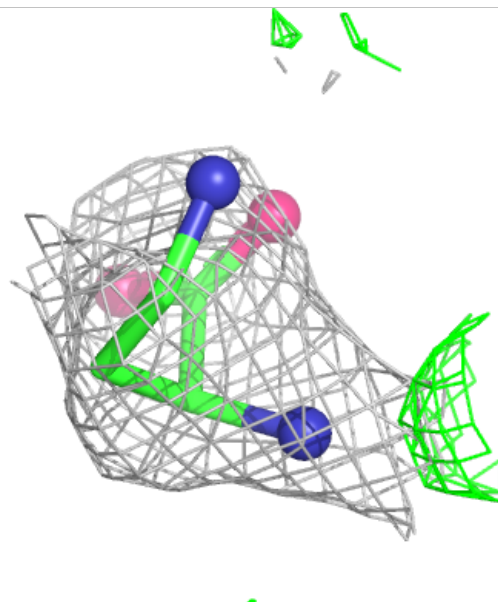
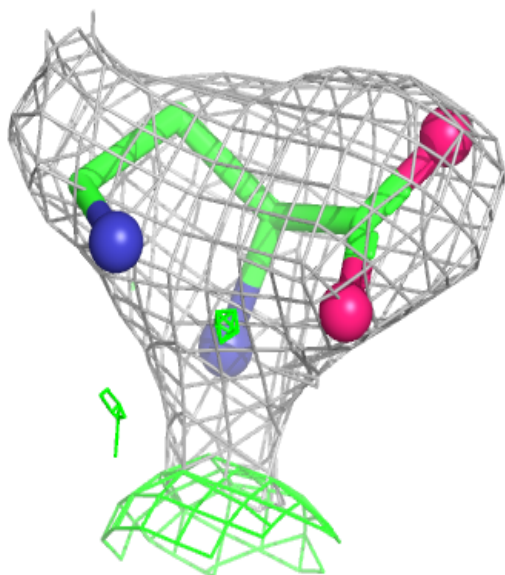
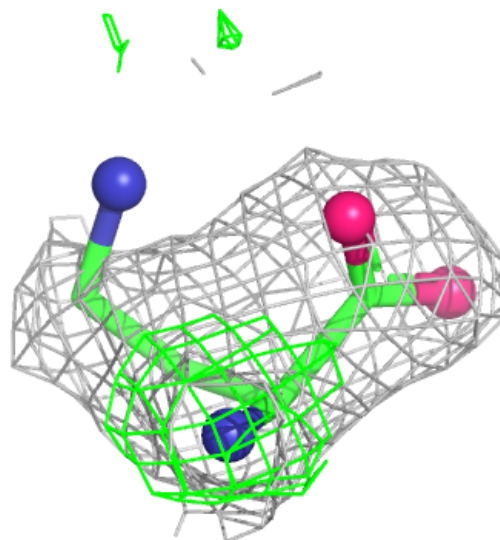
**Electron density around DAB K 401:**

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



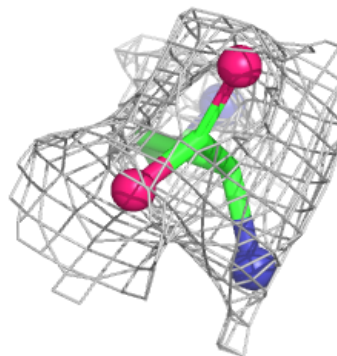
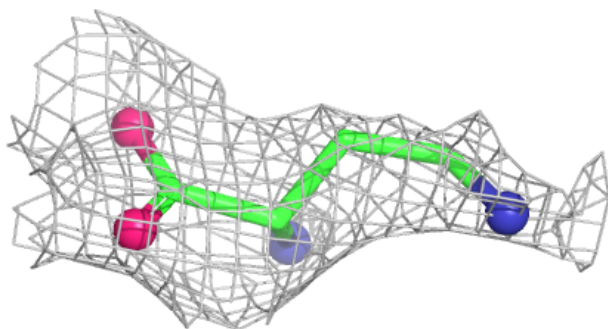
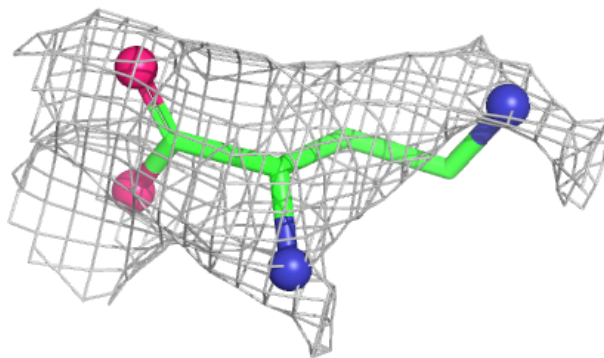
Electron density around DAB F 401:

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



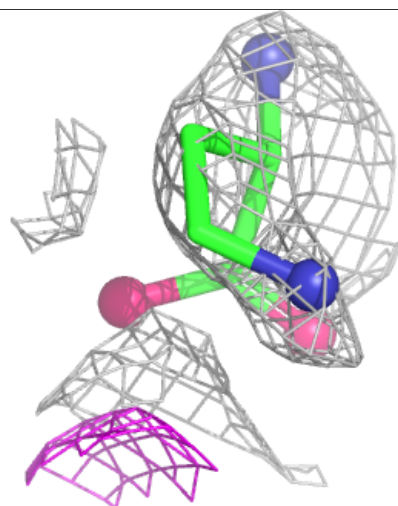
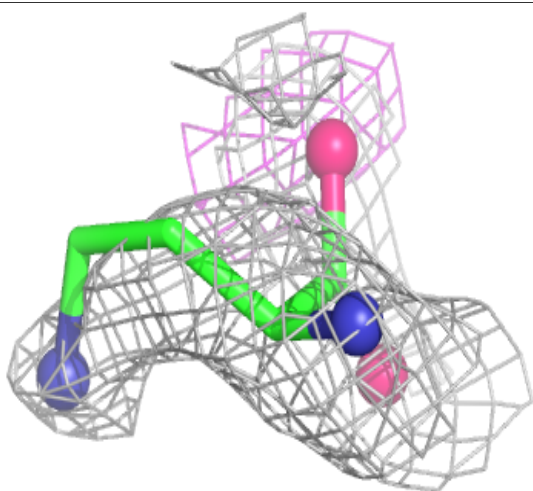
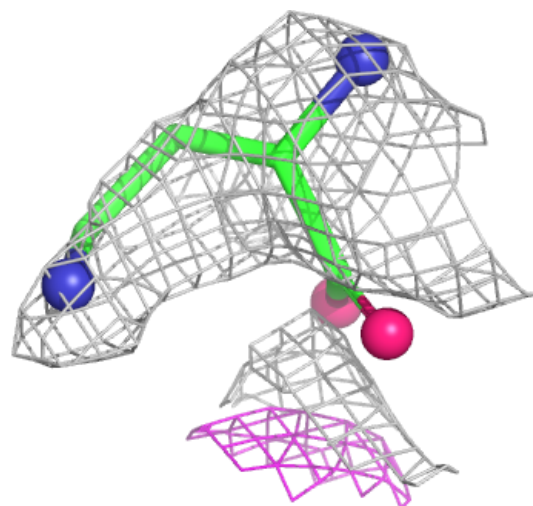
Electron density around DAB A 402:

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



Electron density around DAB J 402:

$2mF_o-DF_c$ (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.