Full wwPDB X-ray Structure Validation Report (i)

Sep 25, 2023 – 07:46 PM EDT

PDB ID	:	6AQJ
Title	:	Crystal structures of Staphylococcus aureus ketol-acid reductoisomerase in complex with two transition state analogs that have biocidal activity.
Authors	:	Patel, K.M.; Teran, D.; Zheng, S.; Gracia, M.; Lv, Y.; Schembri, M.A.; McGeary, R.P.; Schenk, G.; Guddat, L.W.
Deposited on	:	2017-08-20
Resolution	:	1.37 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 (i) 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 (i)) were used in the production of this report:

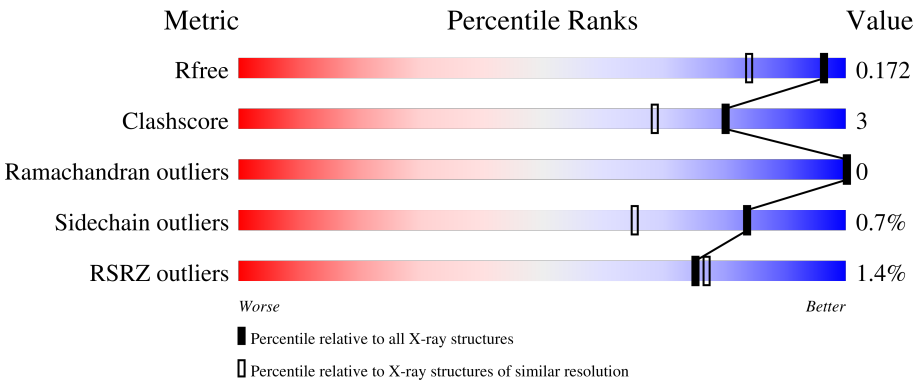
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.37 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}	130704	2907 (1.40-1.36)
Clashscore	141614	3037 (1.40-1.36)
Ramachandran outliers	138981	2970 (1.40-1.36)
Sidechain outliers	138945	2969 (1.40-1.36)
RSRZ outliers	127900	2846 (1.40-1.36)

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

Mol	Chain	Length	Quality of chain
1	A	339	2% 88% 7% •
1	B	339	% 91% 5% •

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 6237 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 Ketol-acid reductoisomerase (NADP(+)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	326	2690	1720	453	509	8	0	23	0
1	B	326	2629	1675	442	501	11	0	12	0

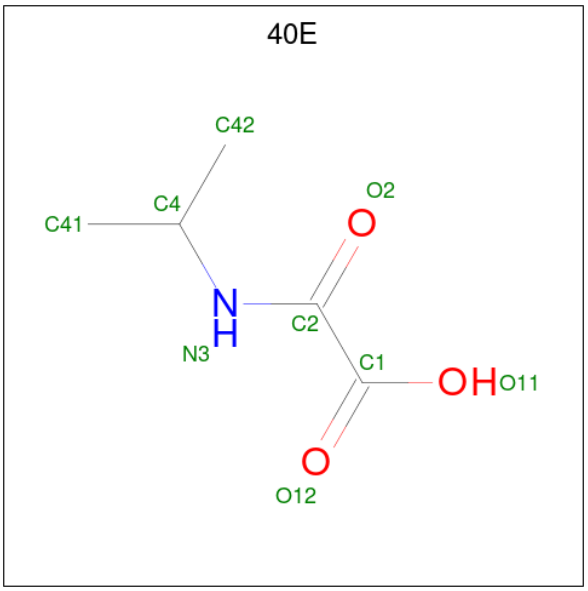
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	335	HIS	-	expression tag	UNP Q2YUF3
A	336	HIS	-	expression tag	UNP Q2YUF3
A	337	HIS	-	expression tag	UNP Q2YUF3
A	338	HIS	-	expression tag	UNP Q2YUF3
A	339	HIS	-	expression tag	UNP Q2YUF3
A	340	HIS	-	expression tag	UNP Q2YUF3
B	335	HIS	-	expression tag	UNP Q2YUF3
B	336	HIS	-	expression tag	UNP Q2YUF3
B	337	HIS	-	expression tag	UNP Q2YUF3
B	338	HIS	-	expression tag	UNP Q2YUF3
B	339	HIS	-	expression tag	UNP Q2YUF3
B	340	HIS	-	expression tag	UNP Q2YUF3

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

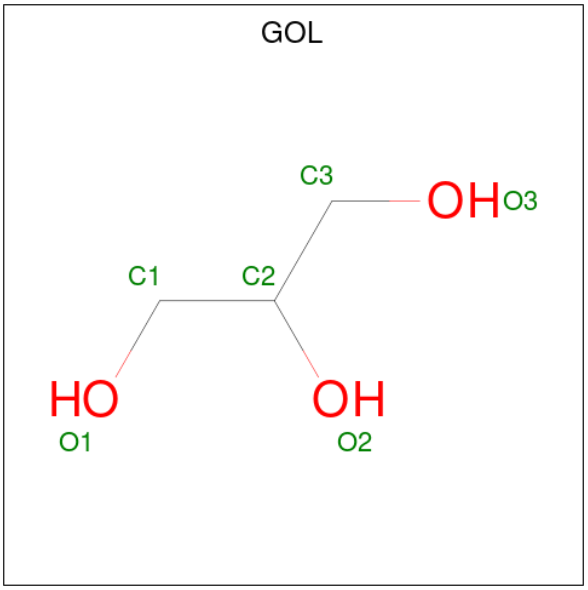
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	2	2	2	0	0
2	B	2	2	2	0	0

- Molecule 3 is oxo(propan-2-ylamino)acetic acid (three-letter code: 40E) (formula: $C_5H_9NO_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	9	5	1	3	0	1
3	B	1	9	5	1	3	0	1

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
4	A	1	6	3	3		0	0
4	A	1	6	3	3		0	0

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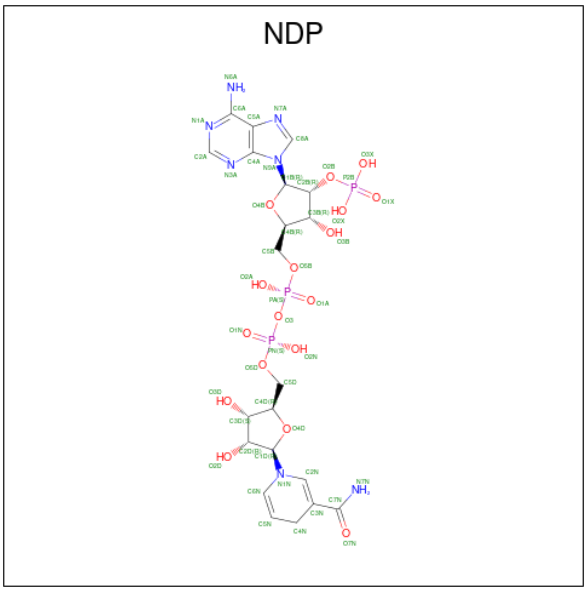
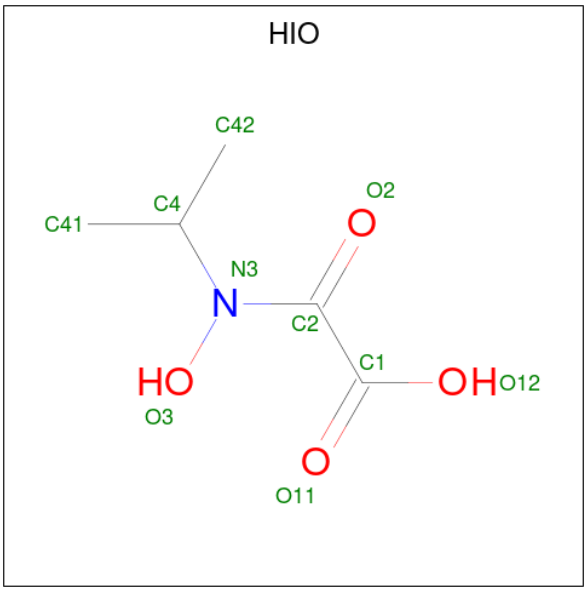
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
4	A	1	6	3	3		0	0
			Total	C	O			
4	B	1	6	3	3		0	0

- Molecule 5 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



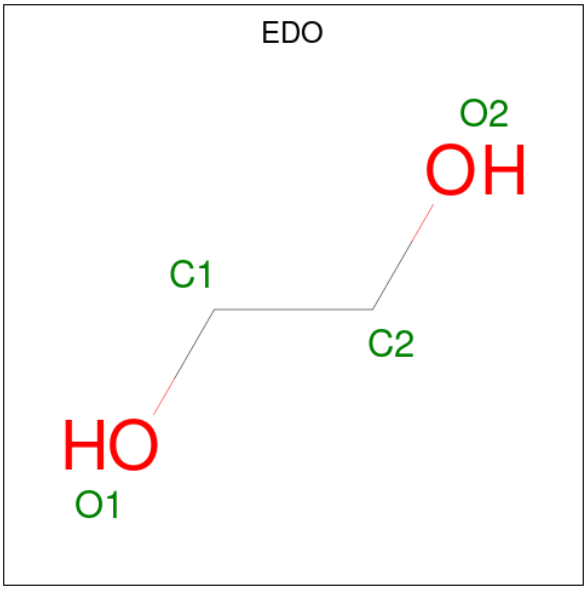
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
5	A	1	48	21	7	17	3	0	0
			Total	C	N	O	P		
5	B	1	48	21	7	17	3	0	0

- Molecule 6 is N-HYDROXY-N-ISOPROPYLOXAMIC ACID (three-letter code: HIO) (formula: $C_5H_9NO_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	A	1	10	5	1	4	0	1
			Total	C	N	O		
6	B	1	10	5	1	4	0	1

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 8 is water.

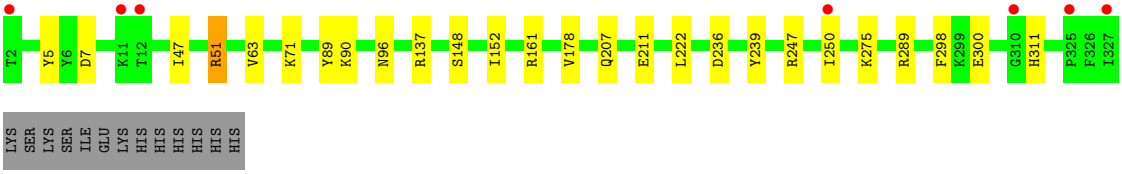
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	362	362	362	0	0
			Total	O		
8	B	386	386	386	0	0

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.

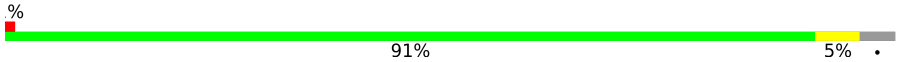
• Molecule 1: Ketol-acid reductoisomerase (NADP(+))

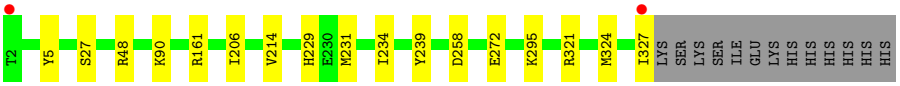
Chain A:



• Molecule 1: Ketol-acid reductoisomerase (NADP(+))

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.09Å 80.73Å 66.64Å 90.00° 92.58° 90.00°	Depositor
Resolution (Å)	34.52 – 1.37 34.52 – 1.37	Depositor EDS
% Data completeness (in resolution range)	97.9 (34.52-1.37) 97.9 (34.52-1.37)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.06 (at 1.37Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.154 , 0.172 0.154 , 0.172	Depositor DCC
R_{free} test set	1991 reflections (1.44%)	wwPDB-VP
Wilson B-factor (Å^2)	13.7	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 51.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.007 for l,k,-h 0.020 for h,-k,-l 0.016 for l,-k,h	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6237	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, NDP, HIO, MG, 40E, GOL

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	0.71	0/2809	0.83	3/3790 (0.1%)
1	B	0.76	1/2715 (0.0%)	0.84	2/3665 (0.1%)
All	All	0.74	1/5524 (0.0%)	0.84	5/7455 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	214	VAL	CB-CG2	-5.24	1.41	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	137	ARG	NE-CZ-NH2	-6.86	116.87	120.30
1	B	48	ARG	NE-CZ-NH1	-6.48	117.06	120.30
1	B	258	ASP	CB-CG-OD2	5.52	123.26	118.30
1	A	236	ASP	CB-CG-OD1	5.43	123.19	118.30
1	A	7	ASP	CB-CG-OD1	5.14	122.93	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2690	0	2699	17	0
1	B	2629	0	2599	9	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	8	1	0
3	B	9	0	8	1	0
4	A	18	0	24	0	0
4	B	6	0	8	0	0
5	A	48	0	26	1	0
5	B	48	0	26	1	0
6	A	10	0	7	1	0
6	B	10	0	7	2	0
7	B	8	0	12	1	0
8	A	362	0	0	1	3
8	B	386	0	0	5	3
All	All	6237	0	5424	29	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (29) 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:275:LYS:HG3	8:B:516:HOH:O	1.74	0.87
1:A:289:ARG:NH2	1:A:300:GLU:OE2	2.18	0.71
1:B:5:TYR:OH	1:B:161[A]:ARG:NH2	2.25	0.69
7:B:406:EDO:H22	8:B:779:HOH:O	1.93	0.69
6:B:402[B]:HIO:C2	5:B:408:NDP:H41N	2.25	0.66
1:B:206:ILE:HD11	1:B:231[B]:MET:SD	2.40	0.62
1:A:247[B]:ARG:NH1	8:A:512:HOH:O	2.37	0.57
1:B:272[A]:GLU:OE2	8:B:503:HOH:O	2.18	0.57
1:A:250:ILE:HG22	6:B:402[B]:HIO:H411	1.87	0.57
1:A:47:ILE:HG22	1:A:63[A]:VAL:HG13	1.90	0.54
1:A:207[B]:GLN:HG3	1:A:311:HIS:NE2	2.23	0.54
1:B:295:LYS:HG3	8:B:606:HOH:O	2.14	0.48
6:A:408[B]:HIO:H411	1:B:234:ILE:HD11	1.97	0.47
1:A:222:LEU:C	1:A:222:LEU:HD23	2.35	0.47
3:B:401[A]:40E:H3	3:B:401[A]:40E:O11	2.15	0.46
1:B:321:ARG:HA	1:B:324[B]:MET:HG3	1.97	0.46
1:A:47:ILE:CG2	1:A:63[A]:VAL:HG13	2.45	0.46
3:A:403[A]:40E:C2	5:A:407:NDP:H41N	2.46	0.46
1:A:51[A]:ARG:HA	1:A:51[A]:ARG:HD2	1.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:TYR:HD2	1:A:90[B]:LYS:HD3	1.82	0.44
1:B:27[A]:SER:OG	8:B:505:HOH:O	2.21	0.44
1:A:239[A]:TYR:CE2	1:B:229:HIS:CE1	3.05	0.44
1:B:324[B]:MET:HB2	1:B:327:ILE:HG12	2.00	0.43
1:A:5:TYR:OH	1:A:161[B]:ARG:NH1	2.52	0.43
1:A:71:LYS:HE2	1:A:96:ASN:OD1	2.20	0.42
1:A:152:ILE:HD11	1:A:161[B]:ARG:HA	2.02	0.42
1:A:152:ILE:HD11	1:A:161[A]:ARG:HA	2.02	0.42
1:A:207[B]:GLN:HG2	1:A:211:GLU:OE2	2.20	0.41
1:A:148:SER:O	1:A:178:VAL:HA	2.20	0.41

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:772:HOH:O	8:B:620:HOH:O[1_556]	2.08	0.12
8:A:762:HOH:O	8:B:778:HOH:O[2_655]	2.12	0.08
8:A:629:HOH:O	8:B:695:HOH:O[2_655]	2.18	0.02

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	348/339 (103%)	340 (98%)	8 (2%)	0	100	100
1	B	337/339 (99%)	329 (98%)	8 (2%)	0	100	100
All	All	685/678 (101%)	669 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	293/282 (104%)	290 (99%)	3 (1%)	76	52
1	B	282/282 (100%)	279 (99%)	3 (1%)	73	48
All	All	575/564 (102%)	569 (99%)	6 (1%)	84	52

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

Mol	Chain	Res	Type
1	A	51[A]	ARG
1	A	51[B]	ARG
1	A	298	PHE
1	B	90	LYS
1	B	239[A]	TYR
1	B	239[B]	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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	EDO	B	407	-	3,3,3	0.49	0	2,2,2	0.68	0
6	HIO	A	408[B]	2	8,9,9	3.05	1 (12%)	8,12,12	2.46	3 (37%)
4	GOL	B	405	-	5,5,5	0.45	0	5,5,5	0.69	0
4	GOL	A	404	-	5,5,5	0.40	0	5,5,5	0.66	0
4	GOL	A	406	-	5,5,5	0.48	0	5,5,5	0.41	0
7	EDO	B	406	-	3,3,3	0.46	0	2,2,2	0.63	0
3	40E	B	401[A]	2	8,8,8	1.57	3 (37%)	9,10,10	2.59	5 (55%)
4	GOL	A	405	-	5,5,5	0.36	0	5,5,5	0.79	0
6	HIO	B	402[B]	2	8,9,9	3.13	1 (12%)	8,12,12	3.09	3 (37%)
5	NDP	B	408	-	45,52,52	1.88	10 (22%)	53,80,80	1.50	9 (16%)
5	NDP	A	407	-	45,52,52	1.26	4 (8%)	53,80,80	1.19	4 (7%)
3	40E	A	403[A]	2	8,8,8	1.79	2 (25%)	9,10,10	1.65	3 (33%)

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	EDO	B	407	-	-	0/1/1/1	-
6	HIO	A	408[B]	2	-	2/8/12/12	-
4	GOL	B	405	-	-	1/4/4/4	-
4	GOL	A	404	-	-	2/4/4/4	-
4	GOL	A	406	-	-	0/4/4/4	-
7	EDO	B	406	-	-	0/1/1/1	-
3	40E	B	401[A]	2	-	2/7/8/8	-
4	GOL	A	405	-	-	0/4/4/4	-
6	HIO	B	402[B]	2	-	4/8/12/12	-
5	NDP	B	408	-	-	5/30/77/77	0/5/5/5
5	NDP	A	407	-	-	4/30/77/77	0/5/5/5
3	40E	A	403[A]	2	-	4/7/8/8	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	402[B]	HIO	C2-N3	8.52	1.43	1.34
6	A	408[B]	HIO	C2-N3	8.36	1.43	1.34
5	B	408	NDP	P2B-O2B	5.43	1.69	1.59
5	B	408	NDP	C5B-C4B	4.68	1.66	1.51
5	A	407	NDP	O4B-C1B	4.23	1.47	1.41
5	B	408	NDP	O5B-C5B	-4.05	1.29	1.44
5	B	408	NDP	C6N-N1N	3.78	1.46	1.37
3	A	403[A]	40E	C2-N3	3.56	1.41	1.34
5	B	408	NDP	C3B-C2B	3.07	1.59	1.52
5	A	407	NDP	P2B-O2B	3.04	1.65	1.59
5	B	408	NDP	C2A-N1A	3.02	1.39	1.33
5	B	408	NDP	C2A-N3A	2.81	1.36	1.32
5	B	408	NDP	O4B-C1B	2.62	1.44	1.41
3	A	403[A]	40E	C2-C1	-2.35	1.51	1.54
3	B	401[A]	40E	C2-N3	2.33	1.38	1.34
5	B	408	NDP	O3B-C3B	2.31	1.48	1.43
3	B	401[A]	40E	C4-N3	-2.27	1.42	1.47
3	B	401[A]	40E	C2-C1	-2.25	1.51	1.54
5	B	408	NDP	C7N-C3N	-2.23	1.43	1.48
5	A	407	NDP	O4D-C1D	2.09	1.46	1.42
5	A	407	NDP	C2A-N3A	2.07	1.35	1.32

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	B	402[B]	HIO	O11-C1-C2	-6.48	114.29	123.07
6	A	408[B]	HIO	C41-C4-N3	-4.68	104.90	109.90
3	B	401[A]	40E	C42-C4-N3	-4.07	101.60	109.79
6	B	402[B]	HIO	O2-C2-N3	-3.82	115.88	120.62
6	B	402[B]	HIO	O12-C1-C2	3.77	122.60	113.84
3	B	401[A]	40E	C4-N3-C2	3.73	127.08	122.74
6	A	408[B]	HIO	O2-C2-N3	-3.58	116.18	120.62
3	B	401[A]	40E	O12-C1-C2	-3.26	114.93	122.18
5	A	407	NDP	C3N-C7N-N7N	3.07	123.13	117.67
5	B	408	NDP	N3A-C2A-N1A	-3.07	123.89	128.68
5	B	408	NDP	C3N-C2N-N1N	-3.03	118.78	123.10
5	B	408	NDP	PN-O3-PA	-2.92	122.81	132.83
5	B	408	NDP	C3B-C2B-C1B	-2.89	97.45	102.89
3	B	401[A]	40E	O11-C1-C2	2.81	121.42	113.15
5	B	408	NDP	O7N-C7N-N7N	-2.78	116.39	122.88
5	B	408	NDP	O2B-P2B-O1X	-2.77	98.68	109.39
5	B	408	NDP	C3N-C7N-N7N	2.77	122.59	117.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403[A]	40E	C4-N3-C2	2.77	125.95	122.74
3	A	403[A]	40E	O2-C2-C1	-2.67	117.56	121.32
3	B	401[A]	40E	O2-C2-N3	-2.61	118.51	123.08
5	B	408	NDP	O5B-PA-O1A	-2.53	99.17	109.07
5	A	407	NDP	C3N-C2N-N1N	-2.31	119.80	123.10
5	B	408	NDP	O4D-C1D-C2D	-2.31	101.60	106.64
6	A	408[B]	HIO	O12-C1-C2	2.28	119.12	113.84
5	A	407	NDP	PN-O3-PA	-2.16	125.42	132.83
3	A	403[A]	40E	C42-C4-N3	-2.06	105.65	109.79
5	A	407	NDP	O7N-C7N-N7N	-2.05	118.08	122.88

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403[A]	40E	O11-C1-C2-N3
3	A	403[A]	40E	C1-C2-N3-C4
3	A	403[A]	40E	O2-C2-N3-C4
3	B	401[A]	40E	C1-C2-N3-C4
3	B	401[A]	40E	O2-C2-N3-C4
5	A	407	NDP	C2B-O2B-P2B-O1X
5	B	408	NDP	C2B-O2B-P2B-O1X
6	A	408[B]	HIO	O11-C1-C2-N3
6	B	402[B]	HIO	C41-C4-N3-O3
6	B	402[B]	HIO	C42-C4-N3-O3
4	A	404	GOL	C1-C2-C3-O3
5	B	408	NDP	C2D-C1D-N1N-C6N
4	A	404	GOL	O2-C2-C3-O3
5	A	407	NDP	C2D-C1D-N1N-C6N
5	B	408	NDP	O4D-C1D-N1N-C6N
6	A	408[B]	HIO	O2-C2-N3-C4
5	A	407	NDP	O4D-C1D-N1N-C6N
4	B	405	GOL	C1-C2-C3-O3
5	B	408	NDP	O4B-C4B-C5B-O5B
3	A	403[A]	40E	O12-C1-C2-N3
5	A	407	NDP	O4B-C4B-C5B-O5B
5	B	408	NDP	C2B-O2B-P2B-O3X
6	B	402[B]	HIO	O11-C1-C2-N3
6	B	402[B]	HIO	C42-C4-N3-C2

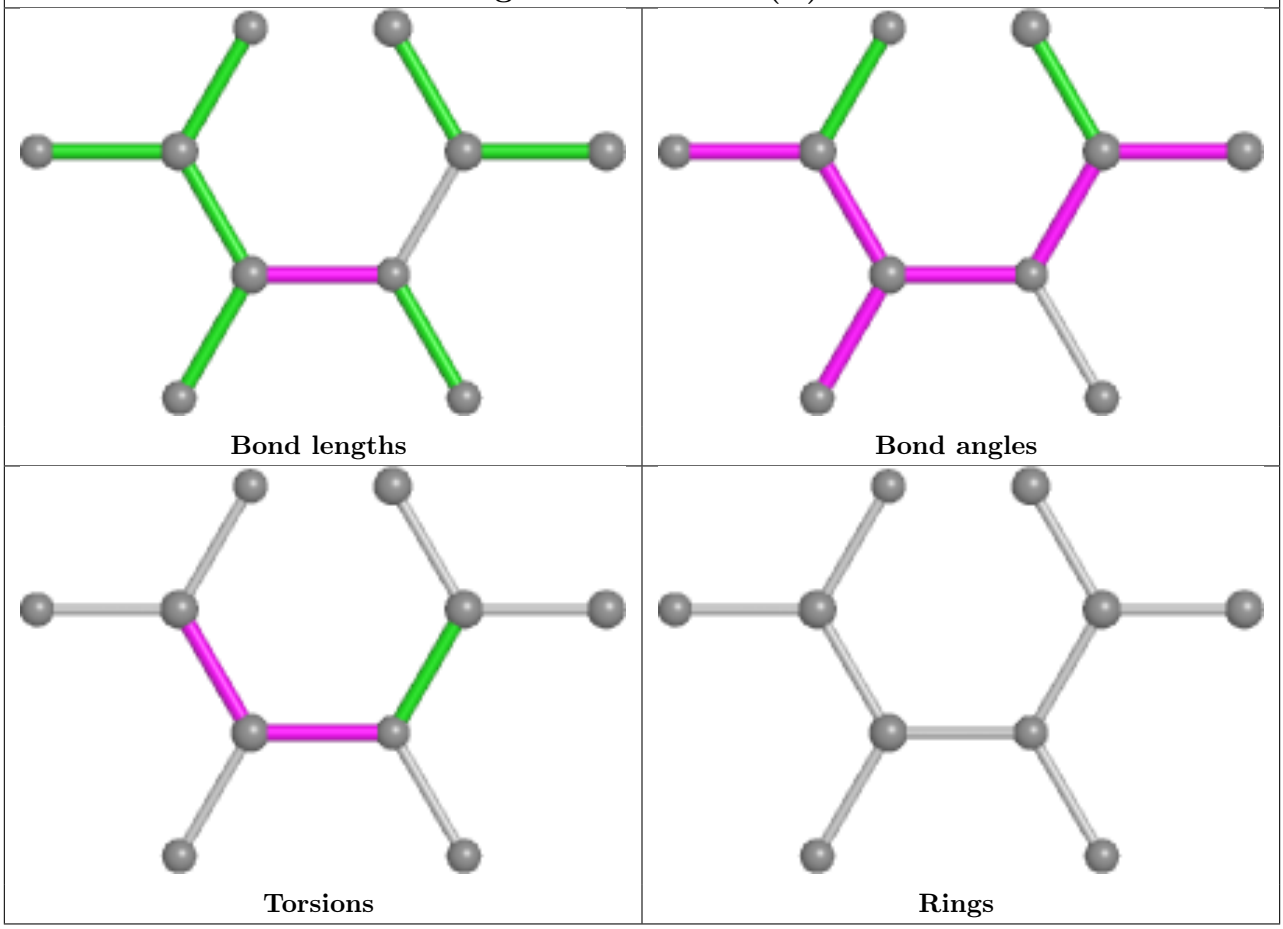
There are no ring outliers.

7 monomers are involved in 6 short contacts:

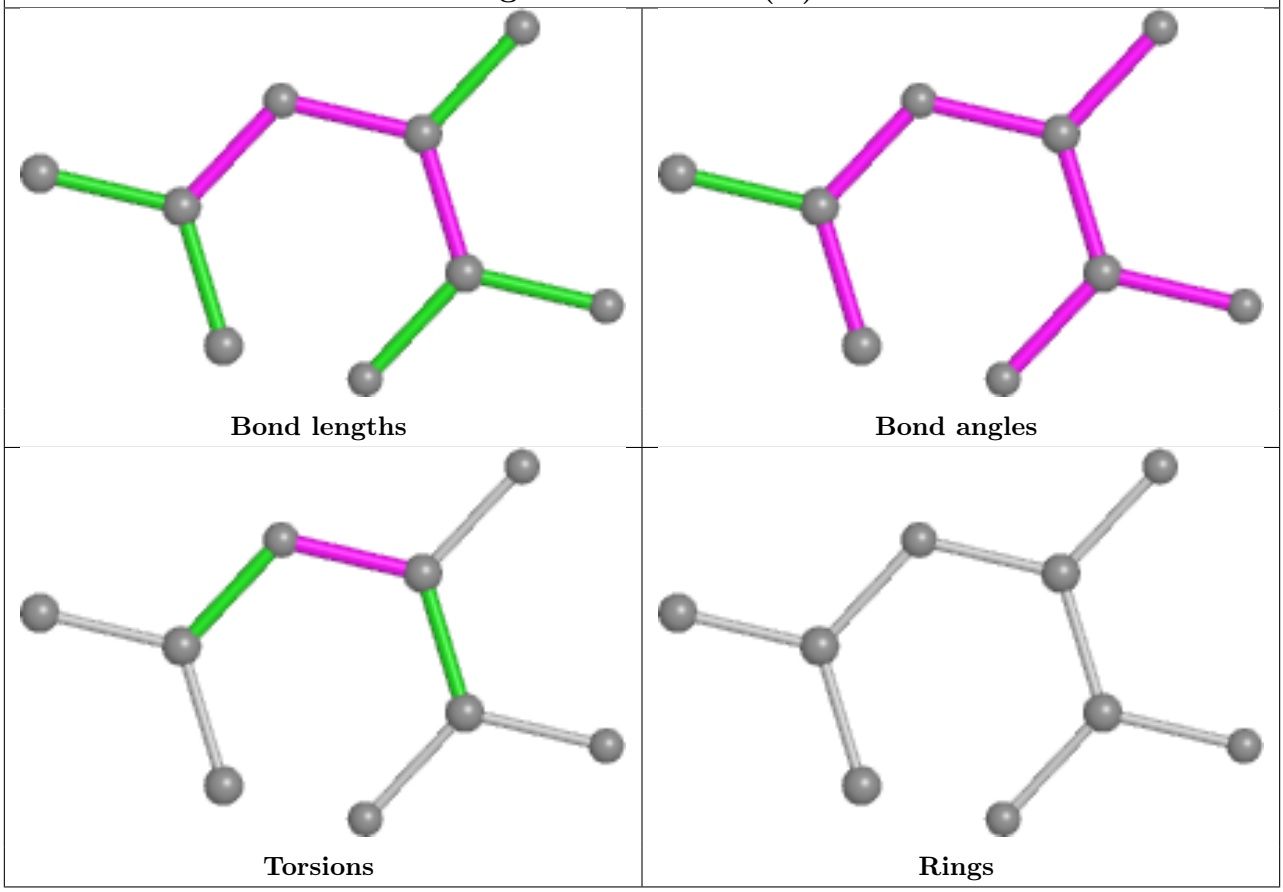
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	408[B]	HIO	1	0
7	B	406	EDO	1	0
3	B	401[A]	40E	1	0
6	B	402[B]	HIO	2	0
5	B	408	NDP	1	0
5	A	407	NDP	1	0
3	A	403[A]	40E	1	0

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

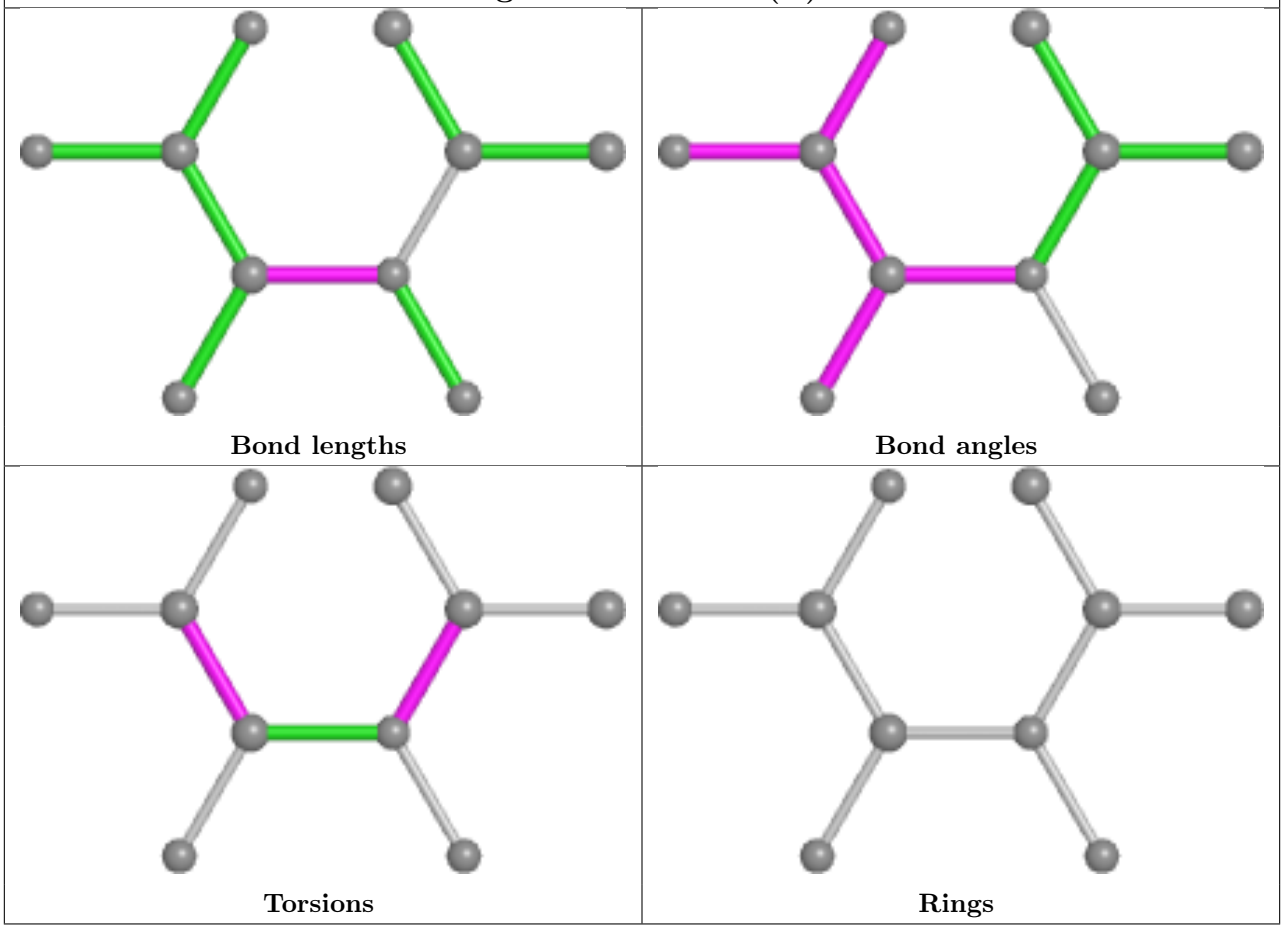
Ligand HIO A 408 (B)



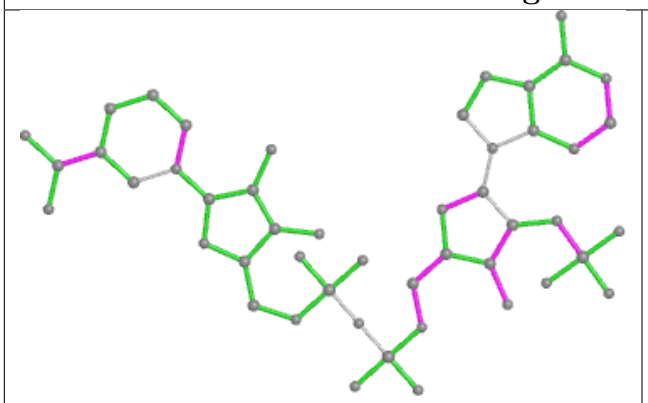
Ligand 40E B 401 (A)



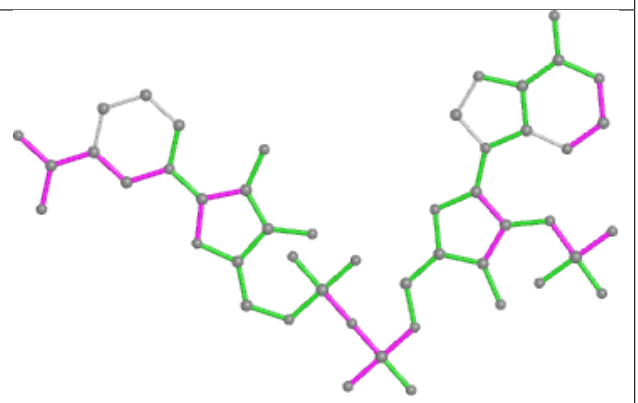
Ligand HIO B 402 (B)



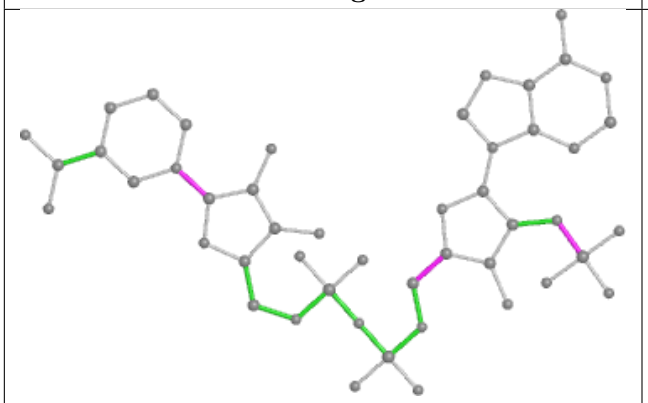
Ligand NDP B 408



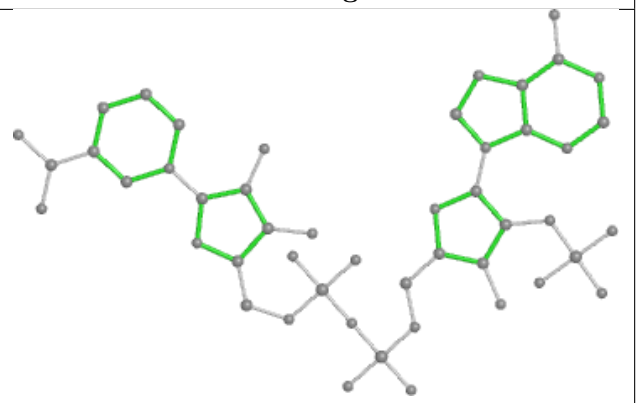
Bond lengths



Bond angles

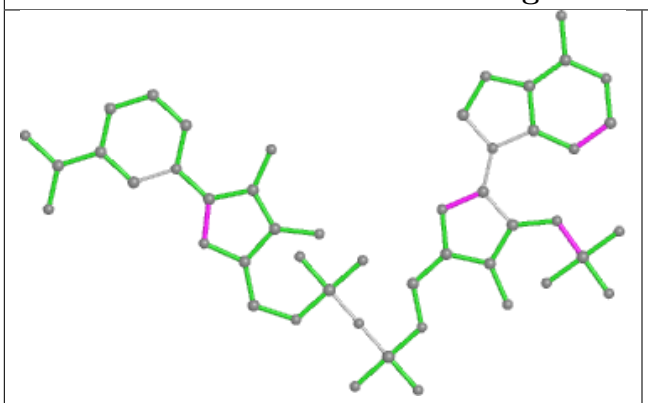


Torsions

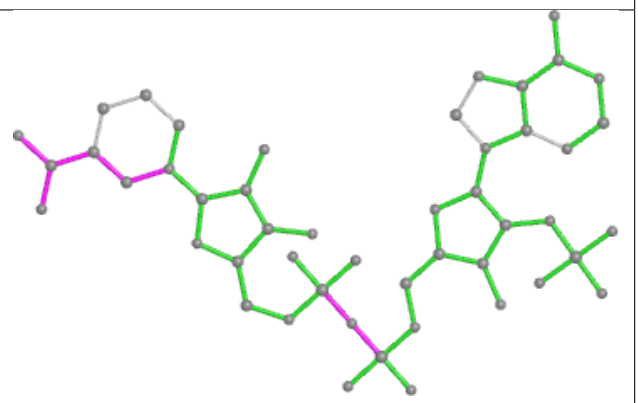


Rings

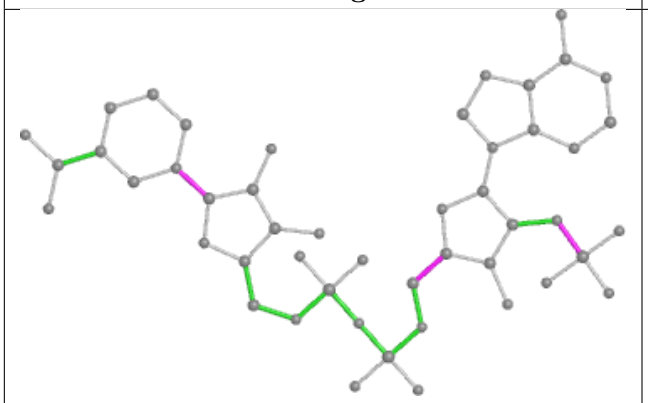
Ligand NDP A 407



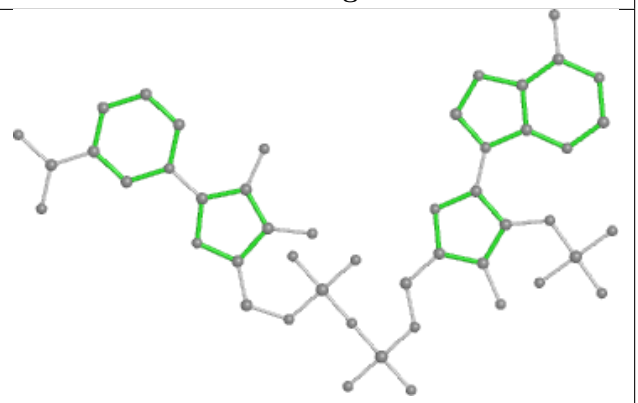
Bond lengths



Bond angles

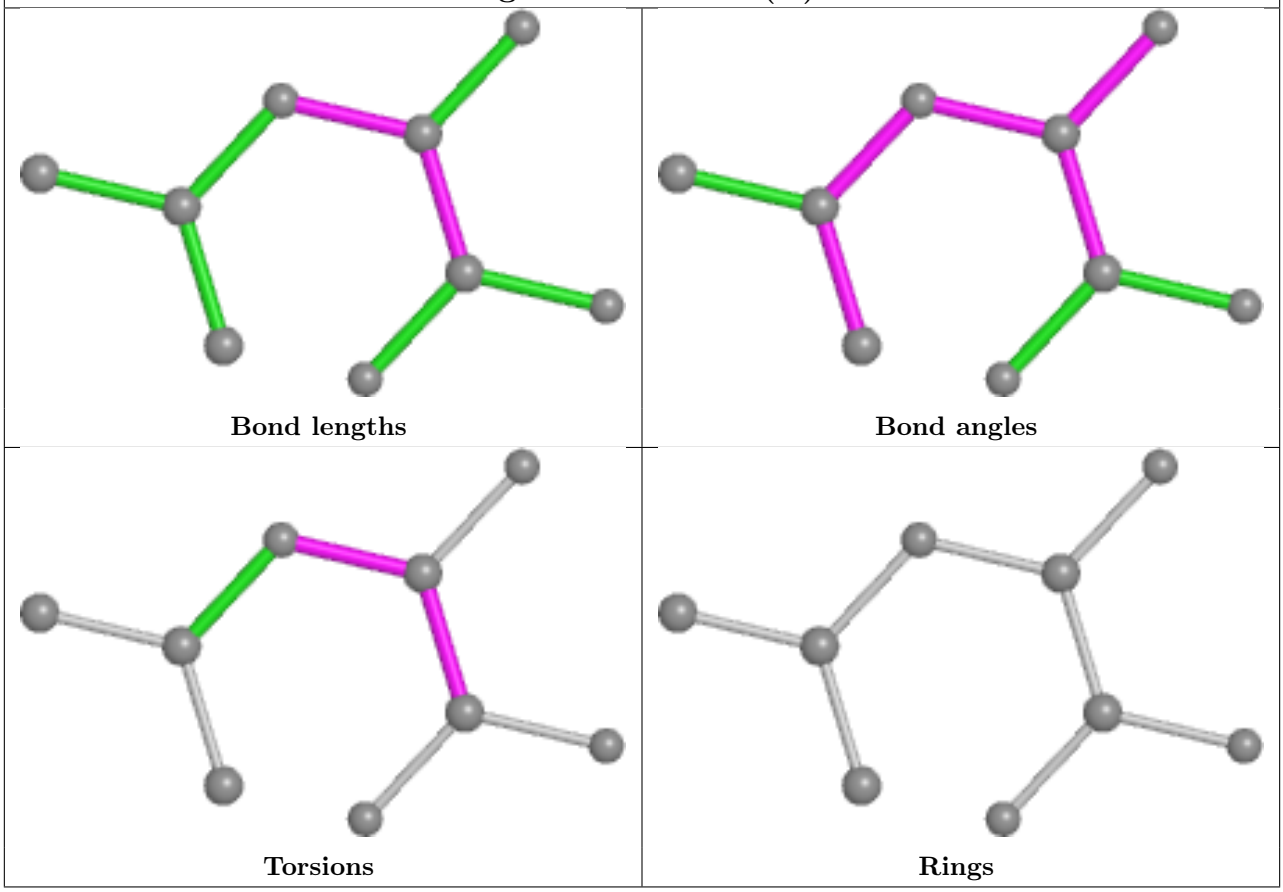


Torsions



Rings

Ligand 40E A 403 (A)



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	326/339 (96%)	0.14	7 (2%)	63	65	9, 16, 30, 49	3 (0%)
1	B	326/339 (96%)	-0.05	2 (0%)	89	90	9, 15, 27, 41	2 (0%)
All	All	652/678 (96%)	0.04	9 (1%)	75	77	9, 15, 29, 49	5 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	THR	4.4
1	A	327	ILE	3.9
1	B	2	THR	3.7
1	B	327	ILE	3.0
1	A	2	THR	2.7
1	A	325	PRO	2.5
1	A	250	ILE	2.3
1	A	11	LYS	2.2
1	A	310	GLY	2.2

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,

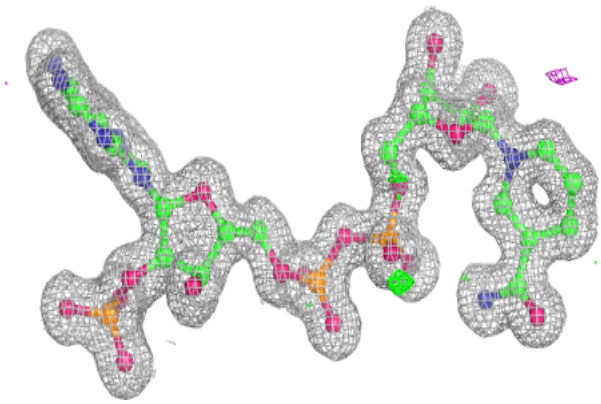
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

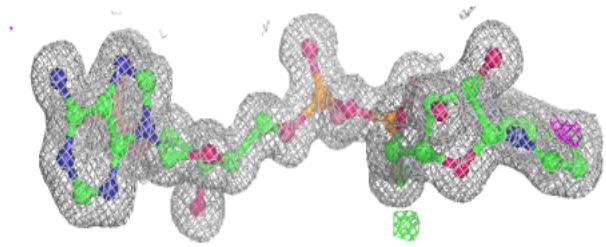
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	EDO	B	407	4/4	0.84	0.12	36,42,46,62	0
7	EDO	B	406	4/4	0.86	0.14	35,39,47,49	0
4	GOL	A	405	6/6	0.89	0.13	23,27,27,32	0
4	GOL	A	404	6/6	0.91	0.13	25,33,36,40	0
4	GOL	B	405	6/6	0.93	0.14	21,27,30,44	0
4	GOL	A	406	6/6	0.97	0.11	19,20,21,24	0
5	NDP	A	407	48/48	0.98	0.07	8,11,16,17	0
5	NDP	B	408	48/48	0.98	0.07	9,12,21,26	0
6	HIO	A	408[B]	10/10	0.98	0.12	7,8,11,13	10
6	HIO	B	402[B]	10/10	0.98	0.12	9,11,15,17	10
3	40E	A	403[A]	9/9	0.98	0.11	7,9,12,13	9
3	40E	B	401[A]	9/9	0.98	0.11	10,10,16,16	9
2	MG	A	401	1/1	0.99	0.04	12,12,12,12	0
2	MG	B	403	1/1	0.99	0.04	12,12,12,12	0
2	MG	A	402	1/1	1.00	0.07	9,9,9,9	0
2	MG	B	404	1/1	1.00	0.08	9,9,9,9	0

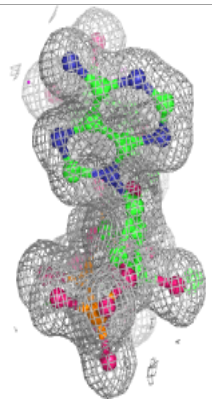
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NDP A 407:

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

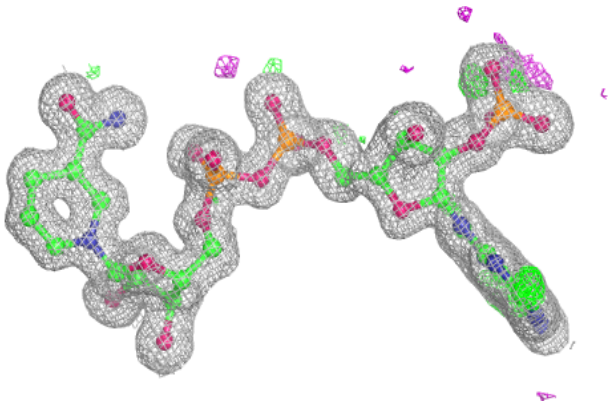


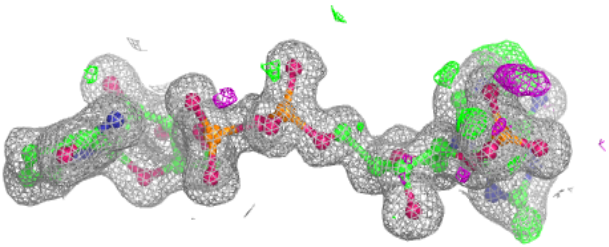


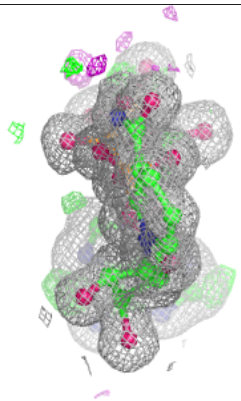


Electron density around NDP B 408:

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

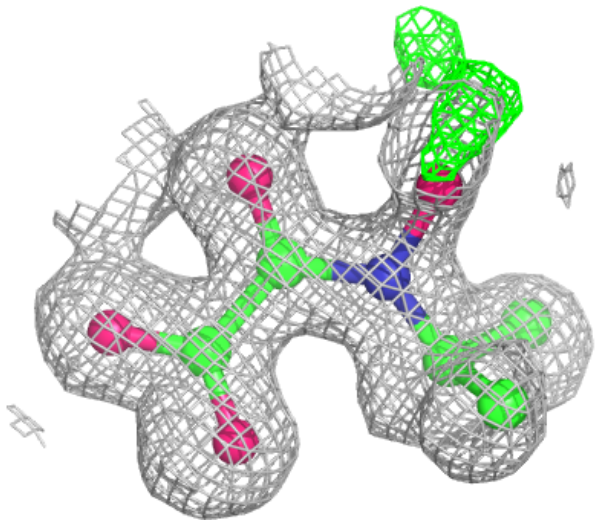


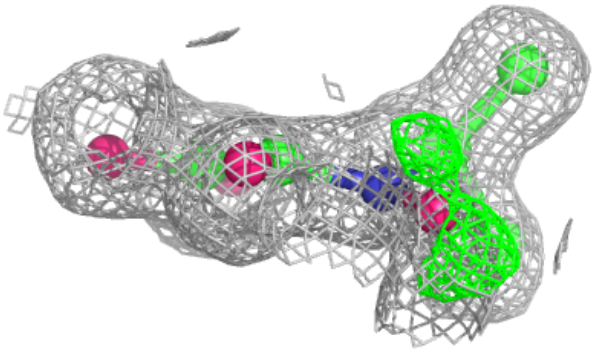


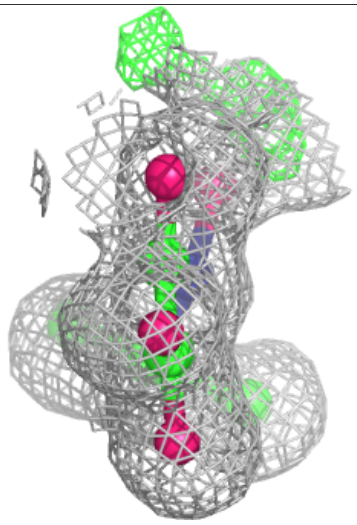


Electron density around HIO A 408 (B):

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







Electron density around HIO B 402 (B):

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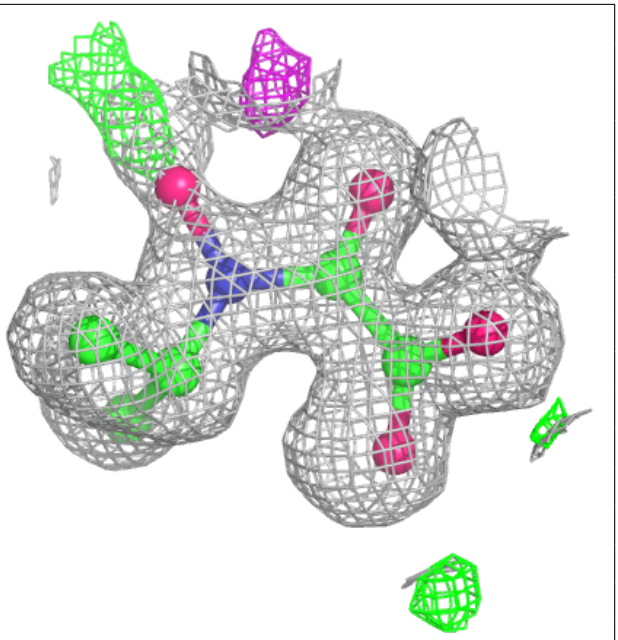
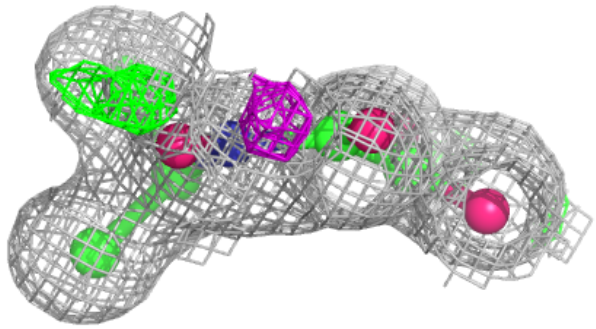
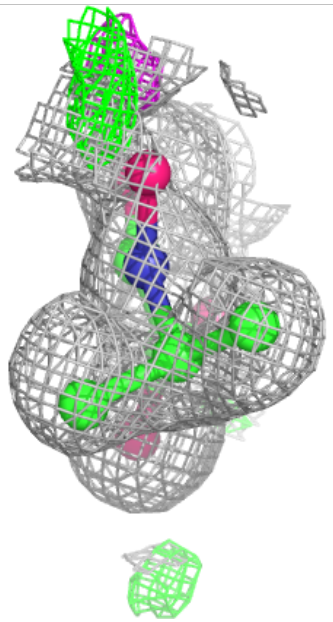
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

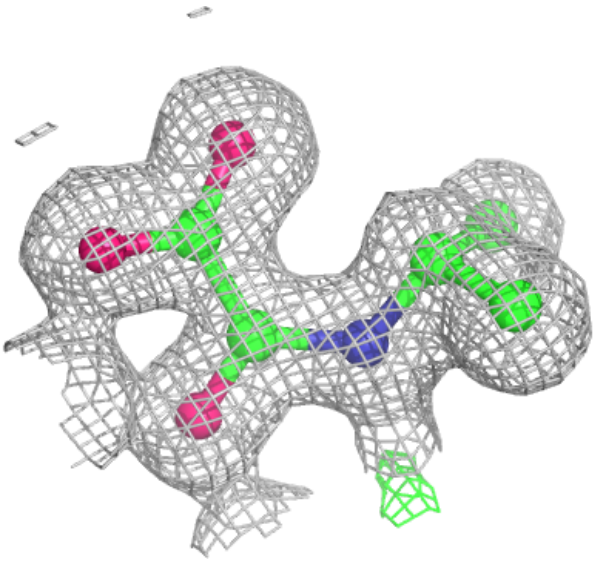


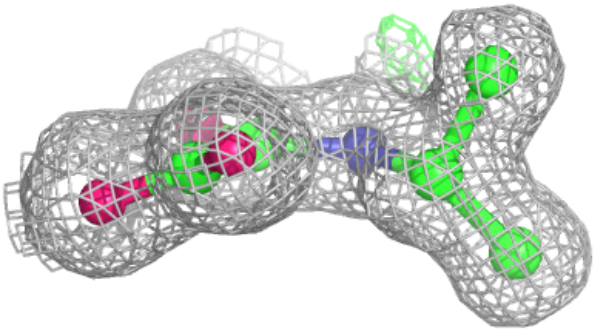


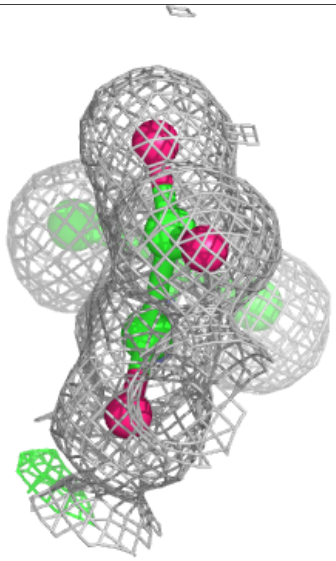


Electron density around 40E A 403 (A):

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

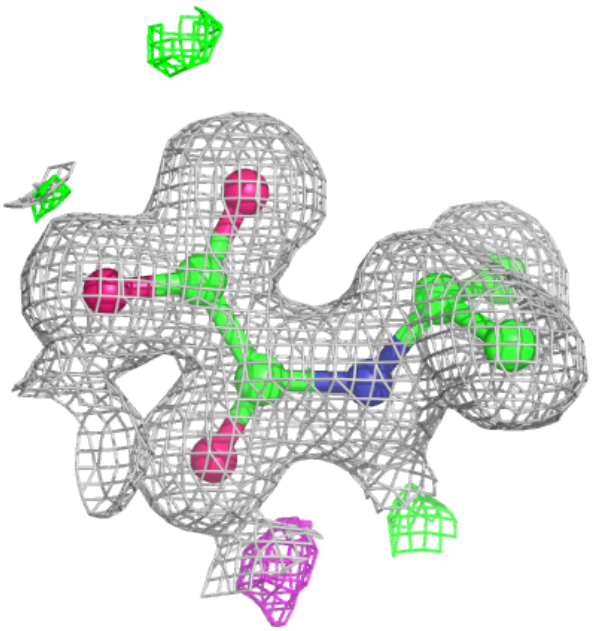


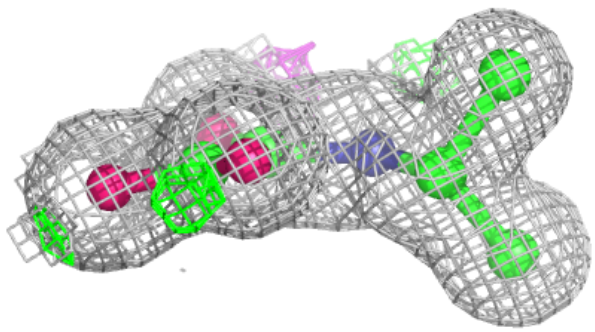


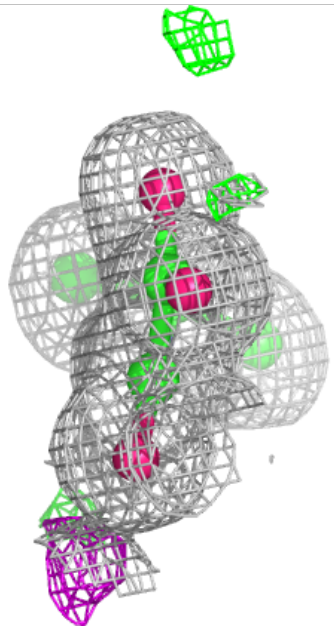


Electron density around 40E B 401 (A):

$2mF_o$-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.