



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:47 PM UTC

PDB ID : 4H80 / pdb_00004h80
Title : Crystal structure of human ALDH3A1 with its isozyme selective inhibitor - N-[4-(4-methylsulfonyl-2-nitroanilino)phenyl]acetamide
Authors : Hurley, T.D.; Parajuli, B.
Deposited on : 2012-09-21
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

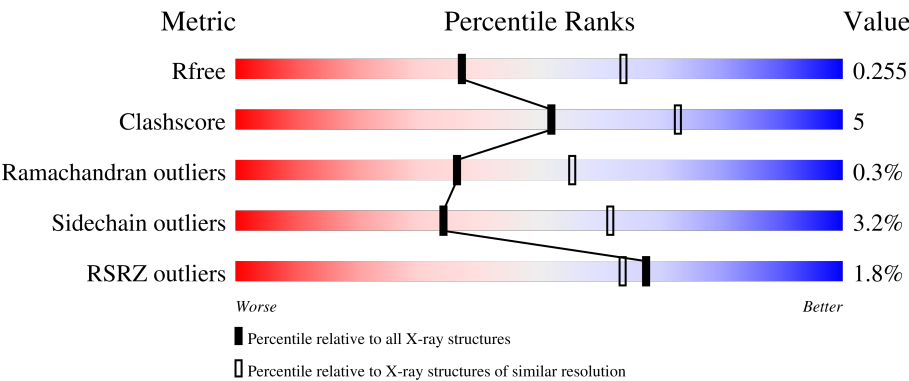
1 Overall quality at a glance i

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	469	3% 85% 9% • 5%
1	B	469	2% 85% 9% • 5%
1	C	469	3% 84% 11% • 5%
1	D	469	% 82% 12% • 5%
1	E	469	% 82% 13% • 5%

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Mol	Chain	Length	Quality of chain
1	F	469	 % 83% 11% 5%
1	G	469	 2% 84% 11% 5%
1	H	469	 % 84% 11% 5%

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
2	04T	A	501	-	-	X	-
2	04T	B	501	-	-	X	-
2	04T	C	501	-	-	X	-
2	04T	F	501	-	-	X	-
2	04T	H	501	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 28414 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 Aldehyde dehydrogenase, dimeric NADP-preferring.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	3	0
			3502	2229	598	657	18			
1	B	446	Total	C	N	O	S	0	3	0
			3504	2231	597	658	18			
1	C	447	Total	C	N	O	S	0	1	0
			3495	2225	595	657	18			
1	D	446	Total	C	N	O	S	0	2	0
			3497	2226	596	657	18			
1	E	446	Total	C	N	O	S	0	3	0
			3506	2232	600	656	18			
1	F	446	Total	C	N	O	S	0	0	0
			3483	2217	592	656	18			
1	G	446	Total	C	N	O	S	0	2	0
			3499	2227	598	656	18			
1	H	447	Total	C	N	O	S	0	4	0
			3517	2239	602	658	18			

There are 136 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	HIS	-	expression tag	UNP P30838
A	-15	HIS	-	expression tag	UNP P30838
A	-14	HIS	-	expression tag	UNP P30838
A	-13	HIS	-	expression tag	UNP P30838
A	-12	HIS	-	expression tag	UNP P30838
A	-11	HIS	-	expression tag	UNP P30838
A	-10	SER	-	expression tag	UNP P30838
A	-9	SER	-	expression tag	UNP P30838
A	-8	GLY	-	expression tag	UNP P30838
A	-7	LEU	-	expression tag	UNP P30838
A	-6	VAL	-	expression tag	UNP P30838
A	-5	PRO	-	expression tag	UNP P30838
A	-4	ARG	-	expression tag	UNP P30838

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P30838
A	-2	SER	-	expression tag	UNP P30838
A	-1	HIS	-	expression tag	UNP P30838
A	133	ALA	SER	engineered mutation	UNP P30838
B	-16	HIS	-	expression tag	UNP P30838
B	-15	HIS	-	expression tag	UNP P30838
B	-14	HIS	-	expression tag	UNP P30838
B	-13	HIS	-	expression tag	UNP P30838
B	-12	HIS	-	expression tag	UNP P30838
B	-11	HIS	-	expression tag	UNP P30838
B	-10	SER	-	expression tag	UNP P30838
B	-9	SER	-	expression tag	UNP P30838
B	-8	GLY	-	expression tag	UNP P30838
B	-7	LEU	-	expression tag	UNP P30838
B	-6	VAL	-	expression tag	UNP P30838
B	-5	PRO	-	expression tag	UNP P30838
B	-4	ARG	-	expression tag	UNP P30838
B	-3	GLY	-	expression tag	UNP P30838
B	-2	SER	-	expression tag	UNP P30838
B	-1	HIS	-	expression tag	UNP P30838
B	133	ALA	SER	engineered mutation	UNP P30838
C	-16	HIS	-	expression tag	UNP P30838
C	-15	HIS	-	expression tag	UNP P30838
C	-14	HIS	-	expression tag	UNP P30838
C	-13	HIS	-	expression tag	UNP P30838
C	-12	HIS	-	expression tag	UNP P30838
C	-11	HIS	-	expression tag	UNP P30838
C	-10	SER	-	expression tag	UNP P30838
C	-9	SER	-	expression tag	UNP P30838
C	-8	GLY	-	expression tag	UNP P30838
C	-7	LEU	-	expression tag	UNP P30838
C	-6	VAL	-	expression tag	UNP P30838
C	-5	PRO	-	expression tag	UNP P30838
C	-4	ARG	-	expression tag	UNP P30838
C	-3	GLY	-	expression tag	UNP P30838
C	-2	SER	-	expression tag	UNP P30838
C	-1	HIS	-	expression tag	UNP P30838
C	133	ALA	SER	engineered mutation	UNP P30838
D	-16	HIS	-	expression tag	UNP P30838
D	-15	HIS	-	expression tag	UNP P30838
D	-14	HIS	-	expression tag	UNP P30838
D	-13	HIS	-	expression tag	UNP P30838

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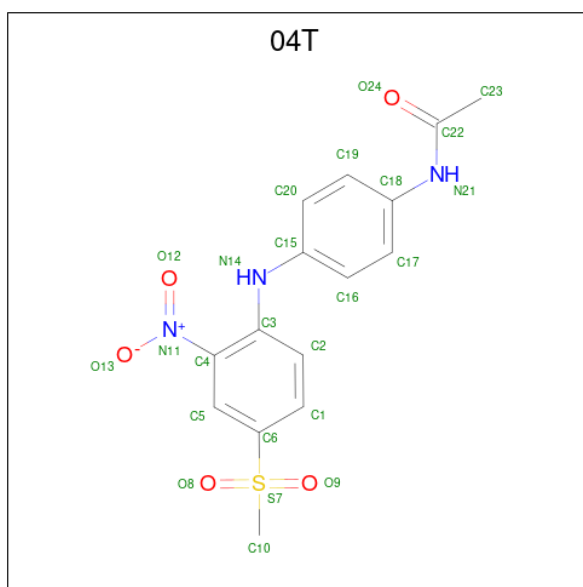
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D	-12	HIS	-	expression tag	UNP P30838
D	-11	HIS	-	expression tag	UNP P30838
D	-10	SER	-	expression tag	UNP P30838
D	-9	SER	-	expression tag	UNP P30838
D	-8	GLY	-	expression tag	UNP P30838
D	-7	LEU	-	expression tag	UNP P30838
D	-6	VAL	-	expression tag	UNP P30838
D	-5	PRO	-	expression tag	UNP P30838
D	-4	ARG	-	expression tag	UNP P30838
D	-3	GLY	-	expression tag	UNP P30838
D	-2	SER	-	expression tag	UNP P30838
D	-1	HIS	-	expression tag	UNP P30838
D	133	ALA	SER	engineered mutation	UNP P30838
E	-16	HIS	-	expression tag	UNP P30838
E	-15	HIS	-	expression tag	UNP P30838
E	-14	HIS	-	expression tag	UNP P30838
E	-13	HIS	-	expression tag	UNP P30838
E	-12	HIS	-	expression tag	UNP P30838
E	-11	HIS	-	expression tag	UNP P30838
E	-10	SER	-	expression tag	UNP P30838
E	-9	SER	-	expression tag	UNP P30838
E	-8	GLY	-	expression tag	UNP P30838
E	-7	LEU	-	expression tag	UNP P30838
E	-6	VAL	-	expression tag	UNP P30838
E	-5	PRO	-	expression tag	UNP P30838
E	-4	ARG	-	expression tag	UNP P30838
E	-3	GLY	-	expression tag	UNP P30838
E	-2	SER	-	expression tag	UNP P30838
E	-1	HIS	-	expression tag	UNP P30838
E	133	ALA	SER	engineered mutation	UNP P30838
F	-16	HIS	-	expression tag	UNP P30838
F	-15	HIS	-	expression tag	UNP P30838
F	-14	HIS	-	expression tag	UNP P30838
F	-13	HIS	-	expression tag	UNP P30838
F	-12	HIS	-	expression tag	UNP P30838
F	-11	HIS	-	expression tag	UNP P30838
F	-10	SER	-	expression tag	UNP P30838
F	-9	SER	-	expression tag	UNP P30838
F	-8	GLY	-	expression tag	UNP P30838
F	-7	LEU	-	expression tag	UNP P30838
F	-6	VAL	-	expression tag	UNP P30838
F	-5	PRO	-	expression tag	UNP P30838

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-4	ARG	-	expression tag	UNP P30838
F	-3	GLY	-	expression tag	UNP P30838
F	-2	SER	-	expression tag	UNP P30838
F	-1	HIS	-	expression tag	UNP P30838
F	133	ALA	SER	engineered mutation	UNP P30838
G	-16	HIS	-	expression tag	UNP P30838
G	-15	HIS	-	expression tag	UNP P30838
G	-14	HIS	-	expression tag	UNP P30838
G	-13	HIS	-	expression tag	UNP P30838
G	-12	HIS	-	expression tag	UNP P30838
G	-11	HIS	-	expression tag	UNP P30838
G	-10	SER	-	expression tag	UNP P30838
G	-9	SER	-	expression tag	UNP P30838
G	-8	GLY	-	expression tag	UNP P30838
G	-7	LEU	-	expression tag	UNP P30838
G	-6	VAL	-	expression tag	UNP P30838
G	-5	PRO	-	expression tag	UNP P30838
G	-4	ARG	-	expression tag	UNP P30838
G	-3	GLY	-	expression tag	UNP P30838
G	-2	SER	-	expression tag	UNP P30838
G	-1	HIS	-	expression tag	UNP P30838
G	133	ALA	SER	engineered mutation	UNP P30838
H	-16	HIS	-	expression tag	UNP P30838
H	-15	HIS	-	expression tag	UNP P30838
H	-14	HIS	-	expression tag	UNP P30838
H	-13	HIS	-	expression tag	UNP P30838
H	-12	HIS	-	expression tag	UNP P30838
H	-11	HIS	-	expression tag	UNP P30838
H	-10	SER	-	expression tag	UNP P30838
H	-9	SER	-	expression tag	UNP P30838
H	-8	GLY	-	expression tag	UNP P30838
H	-7	LEU	-	expression tag	UNP P30838
H	-6	VAL	-	expression tag	UNP P30838
H	-5	PRO	-	expression tag	UNP P30838
H	-4	ARG	-	expression tag	UNP P30838
H	-3	GLY	-	expression tag	UNP P30838
H	-2	SER	-	expression tag	UNP P30838
H	-1	HIS	-	expression tag	UNP P30838
H	133	ALA	SER	engineered mutation	UNP P30838

- Molecule 2 is N-(4-{[4-(methylsulfonyl)-2-nitrophenyl]amino}phenyl)acetamide (CCD ID: 04T) (formula: C₁₅H₁₅N₃O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	B	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	C	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	D	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	E	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	F	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	G	1	Total	C	N	O	S	0	0
			24	15	3	5	1		
2	H	1	Total	C	N	O	S	0	0
			24	15	3	5	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total	O	0	0
			50	50		
3	B	28	Total	O	0	0
			28	28		
3	C	20	Total	O	0	0
			20	20		
3	D	37	Total	O	0	0
			37	37		

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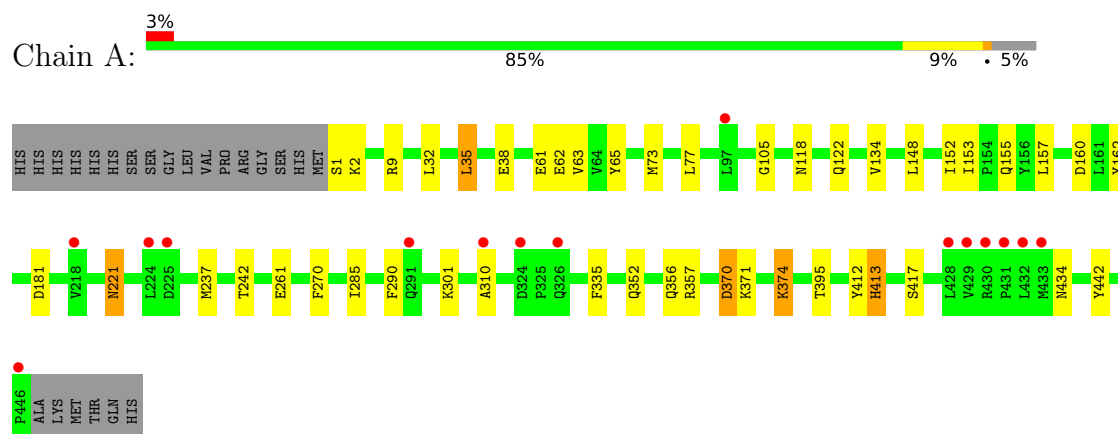
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	32	Total 32	O 32	0	0
3	F	16	Total 16	O 16	0	0
3	G	17	Total 17	O 17	0	0
3	H	19	Total 19	O 19	0	0

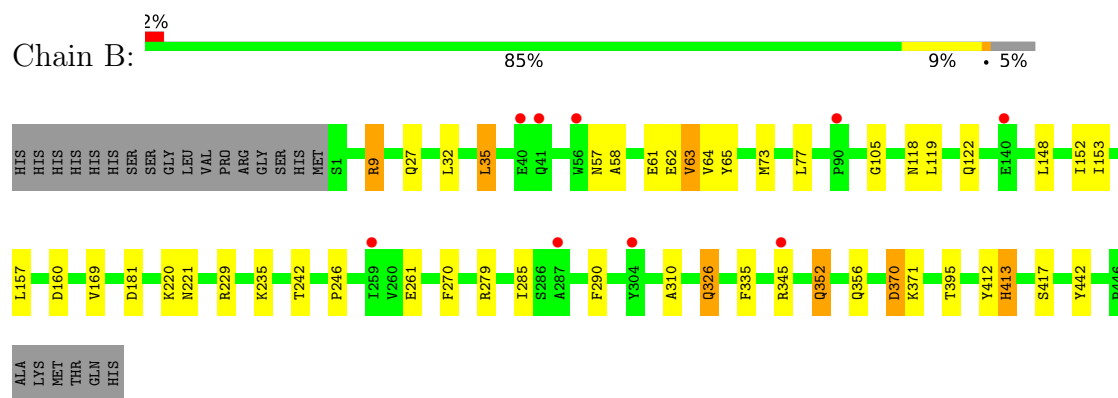
3 Residue-property plots [i](#)

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.

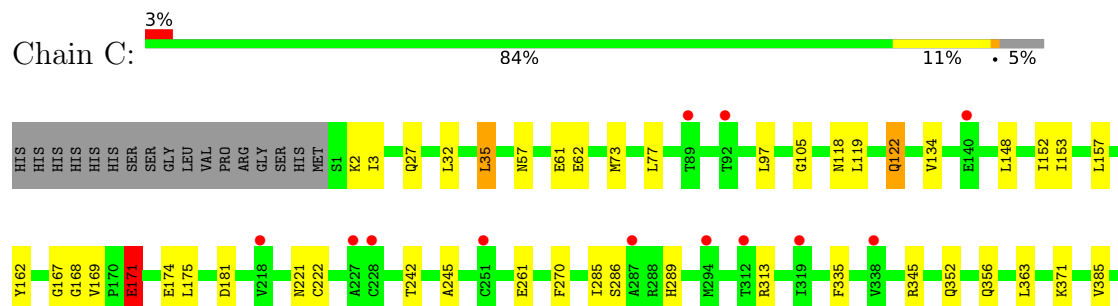
- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring



- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring

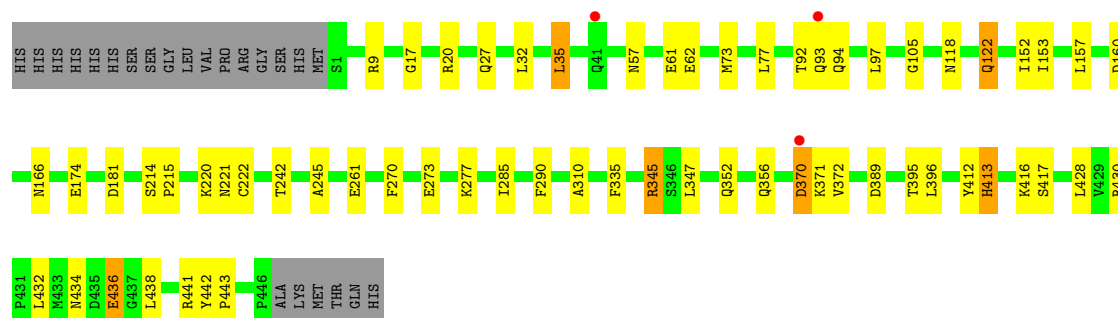
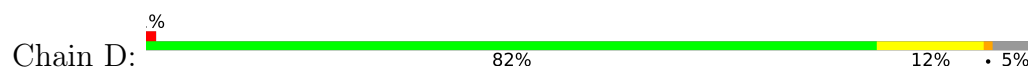


- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring

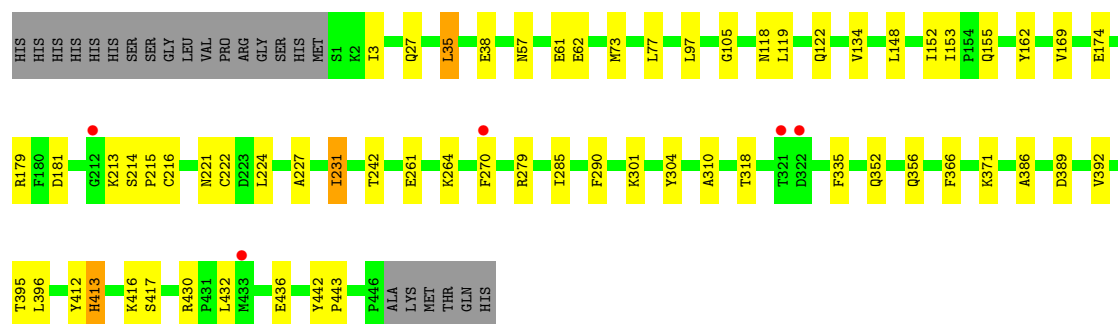
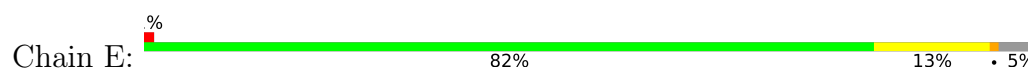




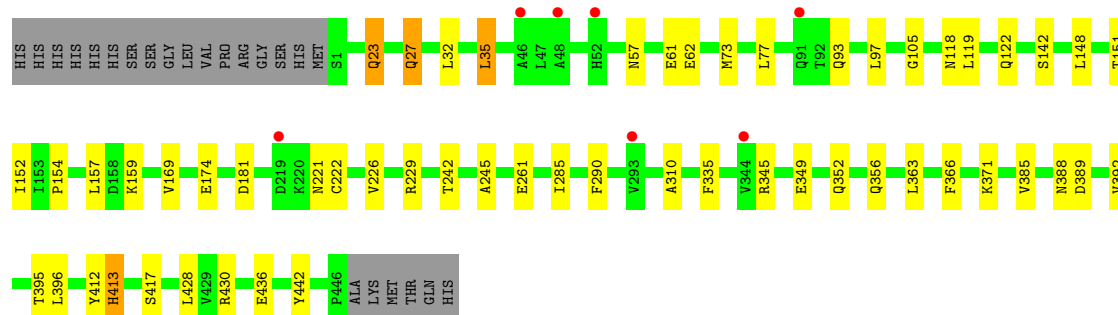
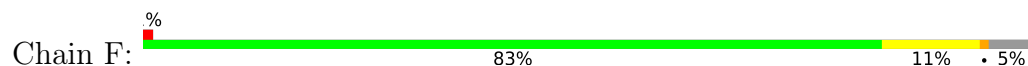
- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring



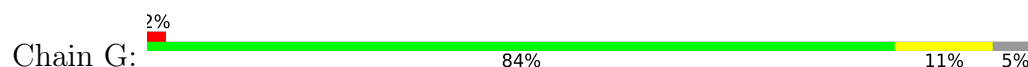
- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring

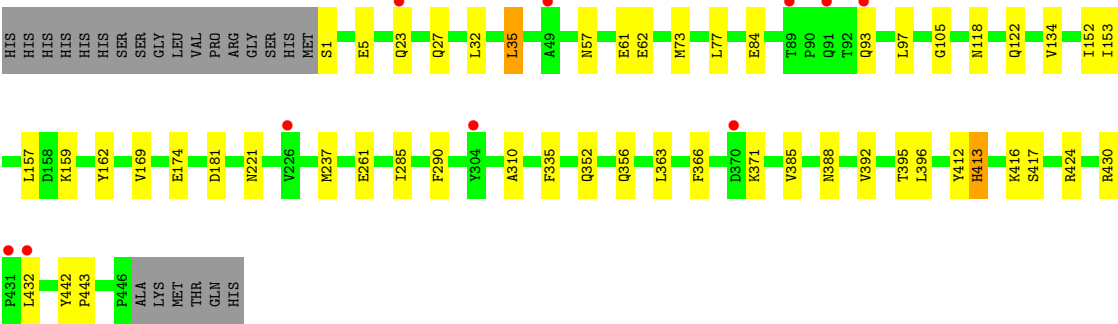


- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring

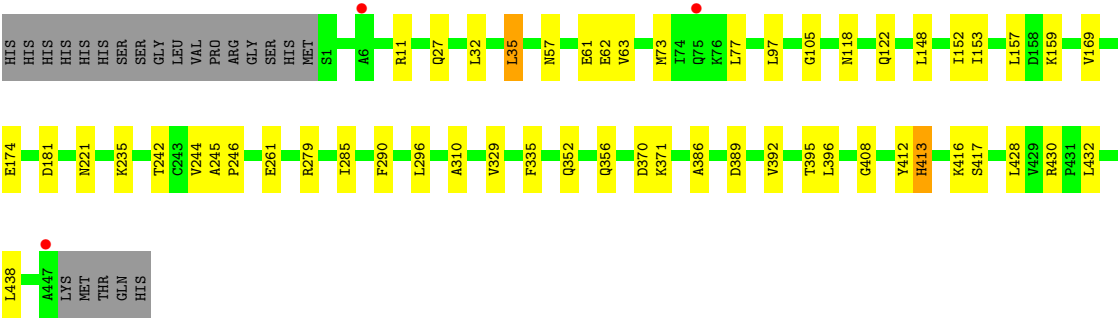
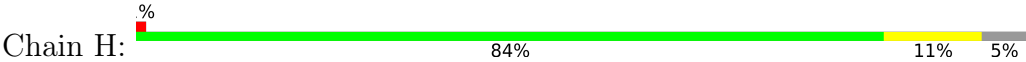


- Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring





● Molecule 1: Aldehyde dehydrogenase, dimeric NADP-preferring



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	89.11Å 95.38Å 117.23Å 112.38° 91.68° 90.99°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.4 (50.00-2.50) 93.5 (50.00-2.50)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.234 , 0.255 0.236 , 0.255	Depositor DCC
R_{free} test set	6195 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	52.5	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.048 for h,-k,-l 0.014 for -h,k,-k-l 0.007 for -h,-k,k+l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	28414	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.88% 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: 04T

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	0.63	2/3584 (0.1%)	0.84	3/4857 (0.1%)
1	B	0.55	4/3587 (0.1%)	0.84	3/4862 (0.1%)
1	C	0.57	4/3572 (0.1%)	0.85	5/4843 (0.1%)
1	D	0.57	3/3576 (0.1%)	0.84	2/4848 (0.0%)
1	E	0.53	3/3589 (0.1%)	0.83	4/4864 (0.1%)
1	F	0.48	1/3556 (0.0%)	0.82	3/4821 (0.1%)
1	G	0.49	2/3578 (0.1%)	0.82	2/4849 (0.0%)
1	H	0.54	2/3603 (0.1%)	0.82	3/4883 (0.1%)
All	All	0.55	21/28645 (0.1%)	0.83	25/38827 (0.1%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	171	GLU	CD-OE1	-11.16	1.04	1.25
1	A	153	ILE	CA-CB	8.50	1.58	1.54
1	G	153	ILE	CA-CB	7.21	1.57	1.54
1	B	352	GLN	CD-NE2	-7.04	1.18	1.33
1	H	153	ILE	CA-CB	6.68	1.57	1.54
1	D	122	GLN	CD-OE1	-6.65	1.10	1.23
1	B	169	VAL	CA-CB	6.54	1.57	1.54
1	D	122	GLN	CD-NE2	-6.28	1.20	1.33
1	B	153	ILE	CA-CB	6.28	1.57	1.54
1	C	153	ILE	CA-CB	6.25	1.57	1.54
1	E	169	VAL	CA-CB	6.12	1.57	1.54
1	D	153	ILE	CA-CB	6.01	1.57	1.54
1	H	169	VAL	CA-CB	5.84	1.57	1.54
1	B	352	GLN	CD-OE1	-5.58	1.12	1.23
1	G	169	VAL	CA-CB	5.54	1.56	1.54
1	E	153	ILE	CA-CB	5.38	1.56	1.54
1	C	169	VAL	CA-CB	5.25	1.56	1.54
1	F	169	VAL	CA-CB	5.24	1.56	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	231	ILE	CG1-CD1	-5.20	1.31	1.51
1	A	63	VAL	C-O	-5.16	1.18	1.24
1	C	122	GLN	C-O	-5.10	1.19	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	171	GLU	CG-CD-OE2	9.64	140.58	118.40
1	C	171	GLU	CG-CD-OE1	-7.13	101.99	118.40
1	E	231	ILE	CB-CG1-CD1	6.80	128.09	113.80
1	E	231	ILE	CB-CA-C	-6.17	104.07	111.97
1	D	416	LYS	CD-CE-NZ	6.08	131.36	111.90
1	C	2	LYS	N-CA-C	-6.06	104.67	111.28
1	G	416	LYS	CD-CE-NZ	5.94	130.91	111.90
1	A	63	VAL	N-CA-C	5.74	116.47	110.62
1	B	63	VAL	N-CA-C	5.57	116.30	110.62
1	C	416	LYS	CD-CE-NZ	5.51	129.53	111.90
1	E	285	ILE	N-CA-C	5.51	115.71	110.42
1	E	231	ILE	CA-CB-CG1	5.44	119.65	110.40
1	F	285	ILE	N-CA-C	5.42	115.62	110.42
1	A	2	LYS	N-CA-C	5.35	117.11	111.28
1	B	326	GLN	CB-CG-CD	5.19	121.42	112.60
1	D	285	ILE	N-CA-C	5.17	115.39	110.42
1	F	442	TYR	CA-C-N	5.17	123.38	119.66
1	F	442	TYR	C-N-CA	5.17	123.38	119.66
1	B	285	ILE	N-CA-C	5.15	115.37	110.42
1	H	63	VAL	N-CA-C	5.11	117.48	111.09
1	A	285	ILE	N-CA-C	5.10	115.32	110.42
1	H	285	ILE	N-CA-C	5.08	115.30	110.42
1	H	408	GLY	N-CA-C	5.05	117.43	111.63
1	C	408	GLY	N-CA-C	5.04	118.12	110.97
1	G	285	ILE	N-CA-C	5.01	115.23	110.42

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3502	0	3539	29	0
1	B	3504	0	3534	25	0
1	C	3495	0	3520	42	0
1	D	3497	0	3529	53	1
1	E	3506	0	3541	45	1
1	F	3483	0	3508	47	0
1	G	3499	0	3534	38	0
1	H	3517	0	3554	42	0
2	A	24	0	15	7	0
2	B	24	0	15	9	0
2	C	24	0	15	9	0
2	D	24	0	15	6	0
2	E	24	0	15	6	0
2	F	24	0	15	8	0
2	G	24	0	15	6	0
2	H	24	0	15	10	0
3	A	50	0	0	5	0
3	B	28	0	0	0	0
3	C	20	0	0	2	0
3	D	37	0	0	2	0
3	E	32	0	0	1	0
3	F	16	0	0	2	0
3	G	17	0	0	1	0
3	H	19	0	0	1	0
All	All	28414	0	28379	287	1

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 5.

All (287) 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:C:167:GLY:HA3	1:C:171:GLU:HG3	1.15	1.09
1:E:179[B]:ARG:HG2	1:E:179[B]:ARG:HH11	1.10	1.07
1:D:20[B]:ARG:HG2	1:D:20[B]:ARG:HH11	1.14	1.06
1:H:279[B]:ARG:HH11	1:H:279[B]:ARG:CG	1.70	1.04
1:H:279[B]:ARG:HH11	1:H:279[B]:ARG:HG2	0.85	1.01
1:E:35:LEU:HD21	1:E:152:ILE:HD12	1.43	1.00
1:H:279[B]:ARG:HG2	1:H:279[B]:ARG:NH1	1.66	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:179[B]:ARG:HH11	1:E:179[B]:ARG:CG	1.76	0.97
1:G:35:LEU:HD21	1:G:152:ILE:HD12	1.47	0.96
1:H:35:LEU:HD21	1:H:152:ILE:HD12	1.47	0.96
1:F:395:THR:HG21	2:F:501:04T:H19	1.45	0.95
1:A:413:HIS:HE1	3:A:626:HOH:O	1.50	0.95
1:D:35:LEU:HD21	1:D:152:ILE:HD12	1.46	0.94
1:F:35:LEU:HD21	1:F:152:ILE:HD12	1.50	0.93
1:C:35:LEU:HD21	1:C:152:ILE:HD12	1.50	0.92
1:A:35:LEU:HD21	1:A:152:ILE:HD12	1.49	0.92
1:B:35:LEU:HD21	1:B:152:ILE:HD12	1.51	0.90
1:C:286:SER:OG	1:C:289:HIS:HB2	1.72	0.88
1:D:20[B]:ARG:HH11	1:D:20[B]:ARG:CG	1.86	0.88
1:G:395:THR:HG21	2:G:501:04T:H19	1.53	0.88
1:F:413:HIS:HE1	3:F:616:HOH:O	1.58	0.87
1:C:242:THR:HG23	2:C:501:04T:H1	1.55	0.87
1:B:9:ARG:HD3	1:B:160:ASP:OD1	1.76	0.85
1:C:167:GLY:CA	1:C:171:GLU:HG3	2.04	0.85
1:C:413:HIS:HD2	3:C:611:HOH:O	1.60	0.84
1:E:224:LEU:HD21	3:E:631:HOH:O	1.77	0.84
1:D:395:THR:HG21	2:D:501:04T:H19	1.60	0.83
1:B:395:THR:HG21	2:B:501:04T:H19	1.60	0.82
1:D:94:GLN:HG3	1:F:151:THR:HG21	1.62	0.80
1:F:97:LEU:HD21	1:H:396:LEU:HD13	1.65	0.78
1:E:179[B]:ARG:HG2	1:E:179[B]:ARG:NH1	1.92	0.77
2:F:501:04T:H19	2:F:501:04T:C23	2.15	0.75
1:B:58:ALA:O	1:B:63:VAL:HG23	1.87	0.75
2:A:501:04T:H19	2:A:501:04T:C23	2.17	0.74
1:E:264:LYS:HD3	1:E:304:TYR:CD1	2.23	0.73
1:F:396:LEU:HD13	1:H:97:LEU:HD21	1.71	0.72
1:D:17:GLY:HA2	1:D:20[B]:ARG:NH1	2.04	0.72
1:B:370:ASP:OD1	1:B:370:ASP:N	2.19	0.72
2:E:501:04T:H19	2:E:501:04T:C23	2.19	0.72
1:D:94:GLN:NE2	1:F:148:LEU:HD12	2.05	0.72
2:D:501:04T:H19	2:D:501:04T:H23A	1.72	0.71
2:H:501:04T:H19	2:H:501:04T:C23	2.21	0.71
2:G:501:04T:H19	2:G:501:04T:C23	2.20	0.71
1:D:97:LEU:HD21	1:E:396:LEU:HD13	1.72	0.71
1:A:412:TYR:O	1:A:413:HIS:HB2	1.91	0.70
2:H:501:04T:H19	2:H:501:04T:H23B	1.73	0.70
1:H:290:PHE:CE1	1:H:310:ALA:HA	2.27	0.70
2:G:501:04T:H19	2:G:501:04T:H23A	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:501:04T:H19	2:B:501:04T:C23	2.22	0.70
1:A:413:HIS:CE1	3:A:626:HOH:O	2.33	0.69
1:B:57:ASN:O	1:B:61:GLU:HB2	1.92	0.69
1:F:290:PHE:CE1	1:F:310:ALA:HA	2.28	0.69
1:G:290:PHE:CE1	1:G:310:ALA:HA	2.28	0.69
1:C:62:GLU:HG3	1:C:118:ASN:HB2	1.74	0.68
1:A:370:ASP:OD1	1:A:370:ASP:N	2.13	0.68
1:F:62:GLU:HG3	1:F:118:ASN:HB2	1.76	0.67
1:A:9[A]:ARG:HD3	1:A:160:ASP:OD1	1.94	0.67
1:D:20[B]:ARG:HG2	1:D:20[B]:ARG:NH1	1.96	0.67
1:G:62:GLU:HG3	1:G:118:ASN:HB2	1.75	0.67
1:D:62:GLU:HG3	1:D:118:ASN:HB2	1.76	0.67
1:D:290:PHE:CE1	1:D:310:ALA:HA	2.29	0.66
1:E:62:GLU:HG3	1:E:118:ASN:HB2	1.76	0.66
1:D:412:TYR:O	1:D:413:HIS:HB2	1.94	0.66
1:A:395:THR:HG21	2:A:501:04T:H19	1.76	0.66
1:F:428:LEU:HD23	1:H:392:VAL:HG23	1.77	0.66
2:D:501:04T:H19	2:D:501:04T:C23	2.25	0.66
1:B:290:PHE:CE1	1:B:310:ALA:HA	2.30	0.65
1:E:290:PHE:CE1	1:E:310:ALA:HA	2.30	0.65
1:D:20[B]:ARG:CG	1:D:20[B]:ARG:NH1	2.51	0.65
1:E:395:THR:HG21	2:E:501:04T:H19	1.77	0.65
1:B:62:GLU:HG3	1:B:118:ASN:HB2	1.78	0.65
1:F:389:ASP:OD2	1:H:430:ARG:NH1	2.31	0.64
1:A:290:PHE:CE1	1:A:310:ALA:HA	2.32	0.63
2:E:501:04T:H19	2:E:501:04T:H23B	1.80	0.63
1:B:412:TYR:O	1:B:413:HIS:HB2	1.98	0.63
2:F:501:04T:H19	2:F:501:04T:H23A	1.80	0.63
1:C:395:THR:HG21	2:C:501:04T:H19	1.81	0.63
1:E:179[B]:ARG:CG	1:E:179[B]:ARG:NH1	2.46	0.62
1:A:242:THR:HG23	2:A:501:04T:H1	1.80	0.62
1:E:412:TYR:O	1:E:413:HIS:HB2	1.99	0.62
1:G:412:TYR:O	1:G:413:HIS:HB2	1.99	0.62
1:D:430:ARG:NH1	1:E:389:ASP:OD2	2.32	0.62
2:C:501:04T:H19	2:C:501:04T:C23	2.30	0.61
1:B:352:GLN:O	1:B:356:GLN:HG3	2.01	0.61
1:C:119:LEU:HD21	2:C:501:04T:H10B	1.81	0.61
1:B:220:LYS:HE3	1:E:38:GLU:HG2	1.83	0.61
1:F:412:TYR:O	1:F:413:HIS:HB2	1.99	0.60
1:F:428:LEU:CD2	1:H:392:VAL:HG23	2.31	0.60
1:D:396:LEU:HD13	1:E:97:LEU:HD21	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:501:04T:H19	2:A:501:04T:H23A	1.83	0.60
2:D:501:04T:H10	3:D:635:HOH:O	2.02	0.60
1:A:270:PHE:HE1	1:B:442:TYR:HB3	1.66	0.60
1:G:413:HIS:HE1	3:G:601:HOH:O	1.85	0.60
1:E:227:ALA:O	1:E:231:ILE:HG12	2.01	0.59
1:C:286:SER:O	1:C:313:ARG:HD2	2.03	0.59
1:D:17:GLY:HA2	1:D:20[B]:ARG:HH12	1.65	0.59
1:D:430:ARG:HD2	1:E:389:ASP:HB2	1.84	0.59
1:D:347:LEU:HD22	1:D:372:VAL:HG22	1.84	0.59
1:H:62:GLU:HG3	1:H:118:ASN:HB2	1.84	0.59
1:H:61:GLU:HG2	2:H:501:04T:H17	1.84	0.58
1:D:94:GLN:HE22	1:F:148:LEU:CD1	2.17	0.58
1:F:389:ASP:HB2	1:H:430:ARG:HD2	1.85	0.58
1:E:231:ILE:HD13	1:E:366:PHE:CE1	2.39	0.58
1:C:168:GLY:H	1:C:171:GLU:HG2	1.68	0.57
1:A:38:GLU:HG2	1:D:220:LYS:HE3	1.87	0.57
1:C:270:PHE:HE1	1:G:442:TYR:HB3	1.69	0.57
1:H:412:TYR:O	1:H:413:HIS:HB2	2.05	0.57
1:F:395:THR:HG21	2:F:501:04T:C19	2.28	0.56
1:C:242:THR:HG23	2:C:501:04T:C1	2.31	0.56
1:B:65:TYR:OH	2:B:501:04T:H23A	2.05	0.56
2:B:501:04T:H19	2:B:501:04T:H23B	1.87	0.55
1:D:389:ASP:OD2	1:E:430:ARG:NH1	2.39	0.55
1:D:270:PHE:HE1	1:E:442:TYR:HB3	1.71	0.55
1:G:352:GLN:O	1:G:356:GLN:HG3	2.06	0.55
1:A:270:PHE:CE1	1:B:442:TYR:HB3	2.42	0.55
1:D:222:CYS:HB2	1:E:432:LEU:HD23	1.88	0.55
1:D:432:LEU:HD23	1:E:222:CYS:HB2	1.88	0.55
1:C:97:LEU:HD21	1:G:396:LEU:HD13	1.88	0.55
1:C:412:TYR:O	1:C:413:HIS:HB2	2.06	0.55
1:D:166:ASN:HA	3:D:623:HOH:O	2.06	0.55
1:F:352:GLN:O	1:F:356:GLN:HG3	2.06	0.55
1:E:264:LYS:HD3	1:E:304:TYR:HD1	1.70	0.54
1:F:222:CYS:HB2	1:H:432:LEU:HD23	1.89	0.54
1:E:352:GLN:O	1:E:356:GLN:HG3	2.08	0.54
1:B:73:MET:HE1	1:B:417:SER:HB2	1.89	0.54
1:E:242:THR:HG23	2:E:501:04T:H1	1.90	0.54
1:H:242:THR:HG23	2:H:501:04T:H1	1.89	0.54
1:E:105:GLY:HA3	1:E:181:ASP:OD1	2.08	0.54
1:H:413:HIS:HE1	3:H:619:HOH:O	1.91	0.54
1:D:94:GLN:HE22	1:F:148:LEU:HD12	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:352:GLN:O	1:C:356:GLN:HG3	2.07	0.53
1:F:345:ARG:HB2	1:G:159:LYS:NZ	2.23	0.53
1:A:155:GLN:OE1	1:D:345:ARG:CB	2.57	0.53
1:A:352:GLN:O	1:A:356:GLN:HG3	2.09	0.53
1:A:62:GLU:HG3	1:A:118:ASN:HB2	1.91	0.53
1:C:396:LEU:HD13	1:G:97:LEU:HD21	1.90	0.52
1:F:345:ARG:CB	1:G:159:LYS:NZ	2.71	0.52
1:G:105:GLY:HA3	1:G:181:ASP:OD1	2.10	0.52
1:A:65:TYR:OH	2:A:501:04T:H23A	2.10	0.52
1:D:389:ASP:HB2	1:E:430:ARG:HD2	1.92	0.52
1:C:105:GLY:HA3	1:C:181:ASP:OD1	2.10	0.52
1:D:436:GLU:HB2	1:F:154:PRO:HG2	1.91	0.52
1:A:155:GLN:OE1	1:D:345:ARG:HB3	2.09	0.52
1:F:23:GLN:HE22	1:F:27:GLN:HE22	1.56	0.52
1:C:222:CYS:HB2	1:G:432:LEU:HD23	1.91	0.51
1:D:352:GLN:O	1:D:356:GLN:HG3	2.10	0.51
1:H:395:THR:HG21	2:H:501:04T:H19	1.92	0.51
1:C:270:PHE:CD1	1:G:443:PRO:HD2	2.45	0.51
2:A:501:04T:H19	2:A:501:04T:H23B	1.92	0.51
1:G:424[A]:ARG:HG2	1:G:424[A]:ARG:HH11	1.75	0.51
1:D:94:GLN:NE2	1:F:148:LEU:CD1	2.72	0.51
1:H:279[B]:ARG:CG	1:H:279[B]:ARG:NH1	2.43	0.51
1:A:442:TYR:HB3	1:B:270:PHE:HE1	1.75	0.51
2:C:501:04T:H19	2:C:501:04T:H23B	1.92	0.51
1:F:73:MET:HE1	1:F:417:SER:HB2	1.93	0.51
1:H:73:MET:HE1	1:H:417:SER:HB2	1.92	0.51
1:H:352:GLN:O	1:H:356:GLN:HG3	2.11	0.51
1:H:57:ASN:O	1:H:61:GLU:HB2	2.10	0.51
1:F:349:GLU:OE1	1:G:159:LYS:HE2	2.10	0.50
1:H:244:VAL:HG22	2:H:501:04T:O9	2.11	0.50
1:H:105:GLY:HA3	1:H:181:ASP:OD1	2.11	0.50
1:G:73:MET:HE1	1:G:417:SER:HB2	1.92	0.50
1:D:105:GLY:HA3	1:D:181:ASP:OD1	2.11	0.50
1:A:73:MET:HE1	1:A:417:SER:HB2	1.93	0.50
1:B:105:GLY:HA3	1:B:181:ASP:OD1	2.12	0.49
2:D:501:04T:C23	2:D:501:04T:C19	2.91	0.49
1:C:395:THR:HG21	2:C:501:04T:H23A	1.95	0.49
1:D:395:THR:HG21	2:D:501:04T:H23A	1.95	0.49
1:G:395:THR:HG21	2:G:501:04T:H23A	1.95	0.49
1:A:105:GLY:HA3	1:A:181:ASP:OD1	2.13	0.49
1:F:413:HIS:CE1	3:F:616:HOH:O	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:LYS:HG3	3:A:641:HOH:O	2.13	0.48
1:D:442:TYR:HB3	1:E:270:PHE:HE1	1.78	0.48
1:B:345[B]:ARG:HB3	1:E:155:GLN:OE1	2.13	0.48
1:C:242:THR:CG2	2:C:501:04T:H1	2.36	0.48
1:A:357:ARG:HD3	3:A:636:HOH:O	2.13	0.47
1:B:395:THR:HG21	2:B:501:04T:C19	2.38	0.47
1:E:73:MET:HE1	1:E:417:SER:HB2	1.96	0.47
2:F:501:04T:H19	2:F:501:04T:H23B	1.92	0.47
1:B:242:THR:HG23	2:B:501:04T:H1	1.97	0.47
1:A:412:TYR:O	1:A:413:HIS:CB	2.62	0.47
1:B:32:LEU:HG	1:B:157:LEU:HD21	1.97	0.47
1:B:119:LEU:HD21	2:B:501:04T:H10B	1.96	0.47
1:C:73:MET:HE1	1:C:417:SER:HB2	1.96	0.47
1:C:389:ASP:OD2	1:G:430:ARG:NH1	2.44	0.47
1:D:443:PRO:HD2	1:E:270:PHE:CD1	2.50	0.46
1:D:73:MET:HE1	1:D:417:SER:HB2	1.96	0.46
2:A:501:04T:C23	2:A:501:04T:C19	2.88	0.46
1:D:428:LEU:CD2	1:E:392:VAL:HG23	2.45	0.46
1:F:392:VAL:HG23	1:H:428:LEU:CD2	2.45	0.46
1:F:119:LEU:HD21	2:F:501:04T:H10B	1.97	0.46
1:C:389:ASP:HB2	1:G:430:ARG:HD2	1.98	0.46
1:D:32:LEU:HG	1:D:157:LEU:HD21	1.98	0.46
1:G:237:MET:HE1	2:G:501:04T:H17	1.97	0.46
1:A:61:GLU:OE2	1:A:237:MET:HE2	2.15	0.45
1:F:226:VAL:O	1:F:229:ARG:HG2	2.16	0.45
1:F:430:ARG:NH1	1:H:389:ASP:OD2	2.46	0.45
1:F:32:LEU:HG	1:F:157:LEU:HD21	1.98	0.45
1:F:345:ARG:HB3	1:G:159:LYS:NZ	2.32	0.45
2:E:501:04T:H19	2:E:501:04T:H23A	1.98	0.45
1:C:168:GLY:H	1:C:171:GLU:CG	2.29	0.45
1:D:428:LEU:O	1:E:386:ALA:HA	2.17	0.45
1:E:148:LEU:O	1:E:152:ILE:HG12	2.16	0.45
1:C:413:HIS:CD2	3:C:611:HOH:O	2.48	0.45
1:C:32:LEU:HG	1:C:157:LEU:HD21	1.98	0.45
1:E:57:ASN:O	1:E:61:GLU:HB2	2.17	0.45
1:G:84:GLU:OE1	1:G:424[B]:ARG:NH2	2.49	0.45
1:E:119:LEU:HD21	2:E:501:04T:H10B	1.98	0.45
1:C:286:SER:OG	1:C:289:HIS:CB	2.55	0.45
1:D:430:ARG:NH1	1:E:392:VAL:HG21	2.31	0.45
1:E:152:ILE:HA	1:E:155:GLN:HE21	1.82	0.44
1:C:168:GLY:N	1:C:171:GLU:HG2	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:363:LEU:HB3	1:C:385:VAL:HG22	1.98	0.44
1:H:242:THR:HG23	2:H:501:04T:C1	2.48	0.44
1:A:221:ASN:OD1	1:A:221:ASN:N	2.51	0.44
1:G:395:THR:HG21	2:G:501:04T:C23	2.47	0.44
1:C:222:CYS:HB2	1:G:432:LEU:CD2	2.47	0.44
1:F:105:GLY:HA3	1:F:181:ASP:OD1	2.18	0.44
1:H:235:LYS:HD3	1:H:246:PRO:O	2.18	0.44
1:D:57:ASN:O	1:D:61:GLU:HB2	2.18	0.44
1:D:270:PHE:CD1	1:E:443:PRO:HD2	2.52	0.44
1:F:363:LEU:HB3	1:F:385:VAL:HG22	1.99	0.43
1:G:424[A]:ARG:HG2	1:G:424[A]:ARG:NH1	2.32	0.43
1:G:93:GLN:OE1	1:G:93:GLN:HA	2.18	0.43
1:A:32:LEU:HG	1:A:157:LEU:HD21	2.00	0.43
1:G:32:LEU:HG	1:G:157:LEU:HD21	2.00	0.43
1:F:57:ASN:O	1:F:61:GLU:HB2	2.18	0.43
1:G:1:SER:O	1:G:5:GLU:HG2	2.17	0.43
1:C:345:ARG:CG	1:H:159:LYS:HZ1	2.31	0.43
1:F:242:THR:HB	1:F:245:ALA:HB2	2.01	0.43
1:H:11[A]:ARG:HG3	1:H:11[A]:ARG:HH11	1.83	0.43
1:G:57:ASN:O	1:G:61:GLU:HB2	2.19	0.43
1:D:214:SER:HA	1:D:215:PRO:HD3	1.89	0.42
1:G:412:TYR:O	1:G:413:HIS:CB	2.67	0.42
2:B:501:04T:C23	2:B:501:04T:C19	2.92	0.42
1:D:9:ARG:CZ	1:D:160:ASP:OD1	2.67	0.42
1:C:345:ARG:HB2	1:H:159:LYS:NZ	2.34	0.42
1:F:428:LEU:O	1:H:386:ALA:HA	2.20	0.42
1:B:148:LEU:O	1:B:152:ILE:HG12	2.18	0.42
1:D:438:LEU:O	1:D:441:ARG:HG2	2.19	0.42
1:C:57:ASN:O	1:C:61:GLU:HB2	2.20	0.42
1:C:285:ILE:HG23	1:C:286:SER:N	2.34	0.42
2:F:501:04T:C23	2:F:501:04T:C19	2.86	0.42
1:C:270:PHE:CE1	1:G:442:TYR:HB3	2.51	0.42
1:D:412:TYR:O	1:D:413:HIS:CB	2.64	0.42
2:C:501:04T:C23	2:C:501:04T:C19	2.97	0.42
1:E:304:TYR:O	1:E:318:THR:HA	2.20	0.42
2:B:501:04T:H19	2:B:501:04T:H23A	1.99	0.42
1:A:148:LEU:O	1:A:152:ILE:HG12	2.20	0.41
1:E:134:VAL:O	1:E:162:TYR:HA	2.20	0.41
1:E:216:CYS:SG	1:E:231:ILE:HD12	2.60	0.41
1:F:392:VAL:HG21	1:H:430:ARG:NH1	2.35	0.41
1:G:363:LEU:HB3	1:G:385:VAL:HG22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:392:VAL:HG12	2:H:501:04T:H23	2.03	0.41
2:H:501:04T:C23	2:H:501:04T:C19	2.88	0.41
1:C:148:LEU:O	1:C:152:ILE:HG12	2.21	0.41
1:C:242:THR:HB	1:C:245:ALA:HB2	2.02	0.41
1:C:428:LEU:HD23	1:G:392:VAL:HG23	2.01	0.41
1:E:412:TYR:O	1:E:413:HIS:CB	2.68	0.41
1:F:222:CYS:HB2	1:H:432:LEU:CD2	2.50	0.41
1:H:32:LEU:HG	1:H:157:LEU:HD21	2.02	0.41
1:D:434:ASN:CG	1:F:159:LYS:HE2	2.45	0.41
1:H:148:LEU:O	1:H:152:ILE:HG12	2.21	0.41
1:C:3:ILE:HG23	1:C:175:LEU:HD13	2.02	0.41
1:F:392:VAL:HG23	1:H:428:LEU:HD23	2.01	0.41
1:F:148:LEU:O	1:F:152:ILE:HG12	2.20	0.41
1:F:345:ARG:HB2	1:G:159:LYS:CE	2.50	0.41
1:H:242:THR:HB	1:H:245:ALA:HB2	2.02	0.41
1:D:370:ASP:OD1	1:D:370:ASP:N	2.54	0.41
1:E:214:SER:HA	1:E:215:PRO:HD3	1.93	0.41
1:G:134:VAL:O	1:G:162:TYR:HA	2.20	0.41
1:A:134:VAL:O	1:A:162:TYR:HA	2.20	0.41
1:H:296:LEU:HB3	1:H:329:VAL:HG23	2.02	0.41
1:H:392:VAL:HG12	2:H:501:04T:C23	2.50	0.41
1:D:9:ARG:NE	1:D:160:ASP:OD1	2.54	0.40
1:F:366:PHE:CD2	1:F:388:ASN:HA	2.56	0.40
1:C:134:VAL:O	1:C:162:TYR:HA	2.21	0.40
1:D:242:THR:HB	1:D:245:ALA:HB2	2.04	0.40
1:H:11[A]:ARG:HG3	1:H:11[A]:ARG:NH1	2.37	0.40
1:D:92:THR:HB	1:D:97:LEU:HD11	2.04	0.40
1:D:428:LEU:HD23	1:E:392:VAL:HG23	2.02	0.40
1:F:93:GLN:HA	1:F:93:GLN:OE1	2.22	0.40
1:F:392:VAL:HG12	2:F:501:04T:C23	2.51	0.40
1:A:357:ARG:HB3	3:A:635:HOH:O	2.20	0.40
1:B:235:LYS:HD3	1:B:246:PRO:O	2.22	0.40
1:B:412:TYR:O	1:B:413:HIS:CB	2.67	0.40
1:G:366:PHE:CD2	1:G:388:ASN:HA	2.57	0.40

All (1) 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:D:273:GLU:OE2	1:E:301:LYS:NZ[1_565]	1.60	0.60

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	447/469 (95%)	438 (98%)	7 (2%)	2 (0%)	30	49
1	B	447/469 (95%)	439 (98%)	7 (2%)	1 (0%)	43	63
1	C	446/469 (95%)	434 (97%)	11 (2%)	1 (0%)	43	63
1	D	446/469 (95%)	437 (98%)	8 (2%)	1 (0%)	43	63
1	E	447/469 (95%)	436 (98%)	10 (2%)	1 (0%)	43	63
1	F	444/469 (95%)	434 (98%)	9 (2%)	1 (0%)	43	63
1	G	446/469 (95%)	437 (98%)	8 (2%)	1 (0%)	43	63
1	H	449/469 (96%)	440 (98%)	8 (2%)	1 (0%)	43	63
All	All	3572/3752 (95%)	3495 (98%)	68 (2%)	9 (0%)	36	55

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	413	HIS
1	C	413	HIS
1	B	413	HIS
1	D	413	HIS
1	E	413	HIS
1	F	413	HIS
1	G	413	HIS
1	H	413	HIS
1	A	434	ASN

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	383/400 (96%)	372 (97%)	11 (3%)	37	65
1	B	383/400 (96%)	369 (96%)	14 (4%)	30	57
1	C	381/400 (95%)	371 (97%)	10 (3%)	40	68
1	D	382/400 (96%)	368 (96%)	14 (4%)	30	57
1	E	383/400 (96%)	369 (96%)	14 (4%)	30	57
1	F	380/400 (95%)	368 (97%)	12 (3%)	34	62
1	G	382/400 (96%)	372 (97%)	10 (3%)	40	68
1	H	384/400 (96%)	372 (97%)	12 (3%)	35	62
All	All	3058/3200 (96%)	2961 (97%)	97 (3%)	34	62

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	35	LEU
1	A	77	LEU
1	A	122	GLN
1	A	221	ASN
1	A	261	GLU
1	A	301	LYS
1	A	335	PHE
1	A	370	ASP
1	A	371	LYS
1	A	374	LYS
1	B	9	ARG
1	B	27	GLN
1	B	35	LEU
1	B	64	VAL
1	B	77	LEU
1	B	122	GLN
1	B	221	ASN
1	B	229	ARG
1	B	261	GLU
1	B	279	ARG
1	B	326	GLN
1	B	335	PHE
1	B	370	ASP
1	B	371	LYS

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Mol	Chain	Res	Type
1	C	27	GLN
1	C	35	LEU
1	C	77	LEU
1	C	122	GLN
1	C	171	GLU
1	C	174	GLU
1	C	221	ASN
1	C	261	GLU
1	C	335	PHE
1	C	371	LYS
1	D	27	GLN
1	D	35	LEU
1	D	77	LEU
1	D	93	GLN
1	D	122	GLN
1	D	174	GLU
1	D	221	ASN
1	D	261	GLU
1	D	277	LYS
1	D	335	PHE
1	D	345	ARG
1	D	370	ASP
1	D	371	LYS
1	D	436	GLU
1	E	3	ILE
1	E	27	GLN
1	E	35	LEU
1	E	77	LEU
1	E	122	GLN
1	E	174	GLU
1	E	213	LYS
1	E	221	ASN
1	E	261	GLU
1	E	279	ARG
1	E	335	PHE
1	E	371	LYS
1	E	416	LYS
1	E	436	GLU
1	F	23	GLN
1	F	27	GLN
1	F	35	LEU
1	F	77	LEU

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Mol	Chain	Res	Type
1	F	122	GLN
1	F	142	SER
1	F	174	GLU
1	F	221	ASN
1	F	261	GLU
1	F	335	PHE
1	F	371	LYS
1	F	436	GLU
1	G	23	GLN
1	G	27	GLN
1	G	35	LEU
1	G	77	LEU
1	G	122	GLN
1	G	174	GLU
1	G	221	ASN
1	G	261	GLU
1	G	335	PHE
1	G	371	LYS
1	H	27	GLN
1	H	35	LEU
1	H	77	LEU
1	H	122	GLN
1	H	174	GLU
1	H	221	ASN
1	H	261	GLU
1	H	335	PHE
1	H	370	ASP
1	H	371	LYS
1	H	416	LYS
1	H	438	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	258	GLN
1	A	326	GLN
1	A	413	HIS
1	B	144	ASN
1	B	155	GLN
1	B	258	GLN
1	B	326	GLN

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Mol	Chain	Res	Type
1	B	423	HIS
1	B	434	ASN
1	C	27	GLN
1	C	39	GLN
1	C	122	GLN
1	C	144	ASN
1	C	155	GLN
1	C	258	GLN
1	C	289	HIS
1	C	326	GLN
1	C	413	HIS
1	D	39	GLN
1	D	91	GLN
1	D	93	GLN
1	D	155	GLN
1	D	258	GLN
1	D	326	GLN
1	E	91	GLN
1	E	258	GLN
1	E	326	GLN
1	F	23	GLN
1	F	27	GLN
1	F	144	ASN
1	F	202	HIS
1	F	258	GLN
1	F	326	GLN
1	F	413	HIS
1	G	27	GLN
1	G	155	GLN
1	G	258	GLN
1	G	326	GLN
1	G	413	HIS
1	H	91	GLN
1	H	144	ASN
1	H	258	GLN
1	H	326	GLN
1	H	413	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	04T	G	501	-	25,25,25	3.55	4 (16%)	32,36,36	1.92	6 (18%)
2	04T	D	501	-	25,25,25	3.64	3 (12%)	32,36,36	1.79	4 (12%)
2	04T	F	501	-	25,25,25	3.49	4 (16%)	32,36,36	2.02	7 (21%)
2	04T	A	501	-	25,25,25	3.57	4 (16%)	32,36,36	2.08	7 (21%)
2	04T	B	501	-	25,25,25	3.57	4 (16%)	32,36,36	1.83	6 (18%)
2	04T	E	501	-	25,25,25	3.37	4 (16%)	32,36,36	1.84	6 (18%)
2	04T	H	501	-	25,25,25	3.17	3 (12%)	32,36,36	1.94	9 (28%)
2	04T	C	501	-	25,25,25	3.47	5 (20%)	32,36,36	1.91	4 (12%)

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	04T	G	501	-	-	8/16/18/18	0/2/2/2
2	04T	D	501	-	-	8/16/18/18	0/2/2/2
2	04T	F	501	-	-	8/16/18/18	0/2/2/2
2	04T	A	501	-	-	10/16/18/18	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	04T	B	501	-	-	10/16/18/18	0/2/2/2
2	04T	E	501	-	-	8/16/18/18	0/2/2/2
2	04T	H	501	-	-	8/16/18/18	0/2/2/2
2	04T	C	501	-	-	9/16/18/18	0/2/2/2

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	04T	C6-S7	-17.42	1.59	1.77
2	A	501	04T	C6-S7	-16.97	1.59	1.77
2	G	501	04T	C6-S7	-16.86	1.59	1.77
2	B	501	04T	C6-S7	-16.86	1.59	1.77
2	F	501	04T	C6-S7	-16.44	1.60	1.77
2	C	501	04T	C6-S7	-16.39	1.60	1.77
2	E	501	04T	C6-S7	-15.93	1.60	1.77
2	H	501	04T	C6-S7	-14.86	1.62	1.77
2	B	501	04T	C4-N11	-3.27	1.39	1.45
2	A	501	04T	C4-N11	-3.27	1.39	1.45
2	F	501	04T	C4-N11	-3.25	1.39	1.45
2	H	501	04T	C4-N11	-3.24	1.39	1.45
2	C	501	04T	C4-N11	-3.11	1.40	1.45
2	E	501	04T	C4-N11	-3.09	1.40	1.45
2	G	501	04T	C4-N11	-2.96	1.40	1.45
2	D	501	04T	C4-N11	-2.94	1.40	1.45
2	C	501	04T	C10-S7	-2.53	1.66	1.75
2	F	501	04T	C10-S7	-2.48	1.66	1.75
2	D	501	04T	C10-S7	-2.47	1.66	1.75
2	F	501	04T	C15-N14	-2.46	1.35	1.40
2	B	501	04T	C10-S7	-2.45	1.66	1.75
2	E	501	04T	C10-S7	-2.44	1.66	1.75
2	G	501	04T	C10-S7	-2.44	1.66	1.75
2	C	501	04T	C15-N14	-2.43	1.35	1.40
2	B	501	04T	C15-N14	-2.42	1.35	1.40
2	H	501	04T	C10-S7	-2.35	1.67	1.75
2	A	501	04T	C15-N14	-2.35	1.35	1.40
2	A	501	04T	C10-S7	-2.34	1.67	1.75
2	G	501	04T	C15-N14	-2.20	1.36	1.40
2	E	501	04T	C15-N14	-2.13	1.36	1.40
2	C	501	04T	O13-N11	-2.01	1.21	1.35

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	04T	C10-S7-C6	6.46	111.67	104.55
2	F	501	04T	C10-S7-C6	6.04	111.21	104.55
2	E	501	04T	C10-S7-C6	6.03	111.19	104.55
2	F	501	04T	O9-S7-O8	-5.87	108.04	117.99
2	D	501	04T	C10-S7-C6	5.81	110.95	104.55
2	C	501	04T	C10-S7-C6	5.68	110.80	104.55
2	G	501	04T	O9-S7-O8	-5.31	108.98	117.99
2	C	501	04T	O9-S7-O8	-5.30	109.00	117.99
2	G	501	04T	C10-S7-C6	5.22	110.30	104.55
2	A	501	04T	O9-S7-O8	-5.17	109.22	117.99
2	H	501	04T	O9-S7-O8	-5.11	109.33	117.99
2	B	501	04T	C10-S7-C6	4.96	110.01	104.55
2	H	501	04T	C10-S7-C6	4.89	109.94	104.55
2	E	501	04T	O9-S7-O8	-4.86	109.74	117.99
2	B	501	04T	O9-S7-O8	-4.66	110.08	117.99
2	B	501	04T	C23-C22-N21	4.64	121.92	114.95
2	A	501	04T	C23-C22-N21	4.41	121.58	114.95
2	D	501	04T	O9-S7-O8	-4.33	110.65	117.99
2	G	501	04T	C23-C22-N21	4.24	121.33	114.95
2	C	501	04T	C23-C22-N21	4.10	121.12	114.95
2	F	501	04T	C23-C22-N21	4.07	121.07	114.95
2	G	501	04T	C1-C6-S7	3.93	123.06	119.58
2	H	501	04T	C23-C22-N21	3.90	120.82	114.95
2	E	501	04T	C23-C22-N21	3.90	120.82	114.95
2	D	501	04T	C1-C6-S7	3.70	122.86	119.58
2	D	501	04T	C23-C22-N21	3.54	120.27	114.95
2	A	501	04T	O24-C22-N21	-3.26	118.59	123.06
2	A	501	04T	C1-C6-S7	3.19	122.41	119.58
2	B	501	04T	C1-C6-S7	2.95	122.20	119.58
2	B	501	04T	O24-C22-N21	-2.95	119.02	123.06
2	H	501	04T	C1-C6-S7	2.90	122.15	119.58
2	A	501	04T	O8-S7-C6	-2.84	106.02	108.23
2	G	501	04T	O9-S7-C6	2.66	110.29	108.23
2	F	501	04T	O9-S7-C6	2.63	110.27	108.23
2	A	501	04T	O9-S7-C6	2.57	110.22	108.23
2	F	501	04T	C1-C6-S7	2.51	121.81	119.58
2	H	501	04T	C2-C1-C6	2.51	121.88	119.44
2	H	501	04T	C1-C6-C5	-2.39	117.72	120.68
2	H	501	04T	O9-S7-C6	2.36	110.06	108.23
2	C	501	04T	O24-C22-N21	-2.35	119.84	123.06
2	H	501	04T	O24-C22-N21	-2.25	119.98	123.06
2	F	501	04T	O24-C22-N21	-2.24	119.99	123.06
2	H	501	04T	O8-S7-C6	2.11	109.87	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	501	04T	C4-C3-N14	-2.11	116.67	119.31
2	G	501	04T	O24-C22-C23	-2.10	118.31	122.05
2	E	501	04T	O9-S7-C6	2.08	109.84	108.23
2	E	501	04T	C5-C6-S7	2.06	120.67	119.02
2	B	501	04T	O9-S7-C6	2.03	109.81	108.23
2	E	501	04T	O24-C22-N21	-2.01	120.30	123.06

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	04T	C3-C4-N11-O12
2	A	501	04T	C5-C4-N11-O12
2	B	501	04T	C5-C4-N11-O12
2	H	501	04T	C3-C4-N11-O12
2	H	501	04T	C5-C4-N11-O12
2	A	501	04T	C23-C22-N21-C18
2	B	501	04T	C23-C22-N21-C18
2	D	501	04T	C23-C22-N21-C18
2	F	501	04T	C23-C22-N21-C18
2	H	501	04T	C23-C22-N21-C18
2	C	501	04T	C23-C22-N21-C18
2	E	501	04T	C23-C22-N21-C18
2	A	501	04T	O24-C22-N21-C18
2	B	501	04T	O24-C22-N21-C18
2	C	501	04T	O24-C22-N21-C18
2	D	501	04T	O24-C22-N21-C18
2	E	501	04T	O24-C22-N21-C18
2	F	501	04T	O24-C22-N21-C18
2	G	501	04T	C23-C22-N21-C18
2	G	501	04T	O24-C22-N21-C18
2	H	501	04T	O24-C22-N21-C18
2	A	501	04T	C5-C6-S7-O8
2	C	501	04T	C5-C6-S7-O8
2	D	501	04T	C5-C6-S7-O8
2	E	501	04T	C5-C6-S7-O8
2	A	501	04T	C1-C6-S7-O8
2	B	501	04T	C1-C6-S7-O8
2	B	501	04T	C5-C6-S7-O8
2	D	501	04T	C1-C6-S7-O8
2	E	501	04T	C1-C6-S7-O8
2	H	501	04T	C1-C6-S7-O8

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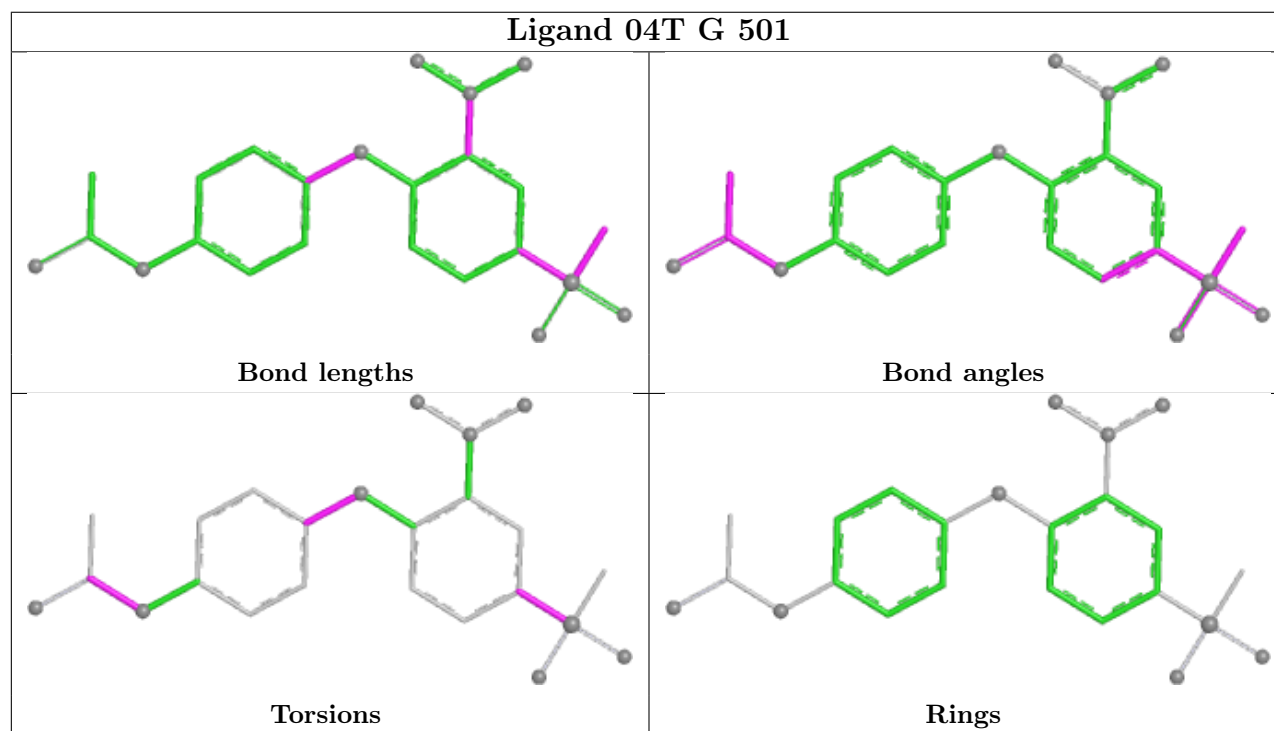
Mol	Chain	Res	Type	Atoms
2	H	501	04T	C5-C6-S7-O8
2	C	501	04T	C1-C6-S7-O8
2	G	501	04T	C5-C6-S7-O8
2	A	501	04T	C5-C6-S7-C10
2	C	501	04T	C5-C6-S7-C10
2	A	501	04T	C1-C6-S7-C10
2	B	501	04T	C5-C6-S7-C10
2	C	501	04T	C1-C6-S7-C10
2	H	501	04T	C5-C6-S7-C10
2	F	501	04T	C5-C6-S7-O8
2	G	501	04T	C1-C6-S7-O8
2	B	501	04T	C1-C6-S7-C10
2	E	501	04T	C5-C6-S7-C10
2	H	501	04T	C1-C6-S7-C10
2	F	501	04T	C1-C6-S7-O8
2	D	501	04T	C1-C6-S7-C10
2	D	501	04T	C5-C6-S7-C10
2	G	501	04T	C16-C15-N14-C3
2	G	501	04T	C20-C15-N14-C3
2	D	501	04T	C16-C15-N14-C3
2	D	501	04T	C20-C15-N14-C3
2	B	501	04T	C16-C15-N14-C3
2	E	501	04T	C1-C6-S7-C10
2	A	501	04T	C16-C15-N14-C3
2	G	501	04T	C5-C6-S7-C10
2	B	501	04T	C20-C15-N14-C3
2	A	501	04T	C20-C15-N14-C3
2	G	501	04T	C1-C6-S7-C10
2	C	501	04T	C5-C4-N11-O12
2	E	501	04T	C5-C4-N11-O12
2	F	501	04T	C16-C15-N14-C3
2	F	501	04T	C20-C15-N14-C3
2	B	501	04T	C3-C4-N11-O12
2	C	501	04T	C3-C4-N11-O12
2	E	501	04T	C3-C4-N11-O12
2	F	501	04T	C1-C6-S7-C10
2	C	501	04T	C4-C3-N14-C15
2	F	501	04T	C5-C6-S7-C10

There are no ring outliers.

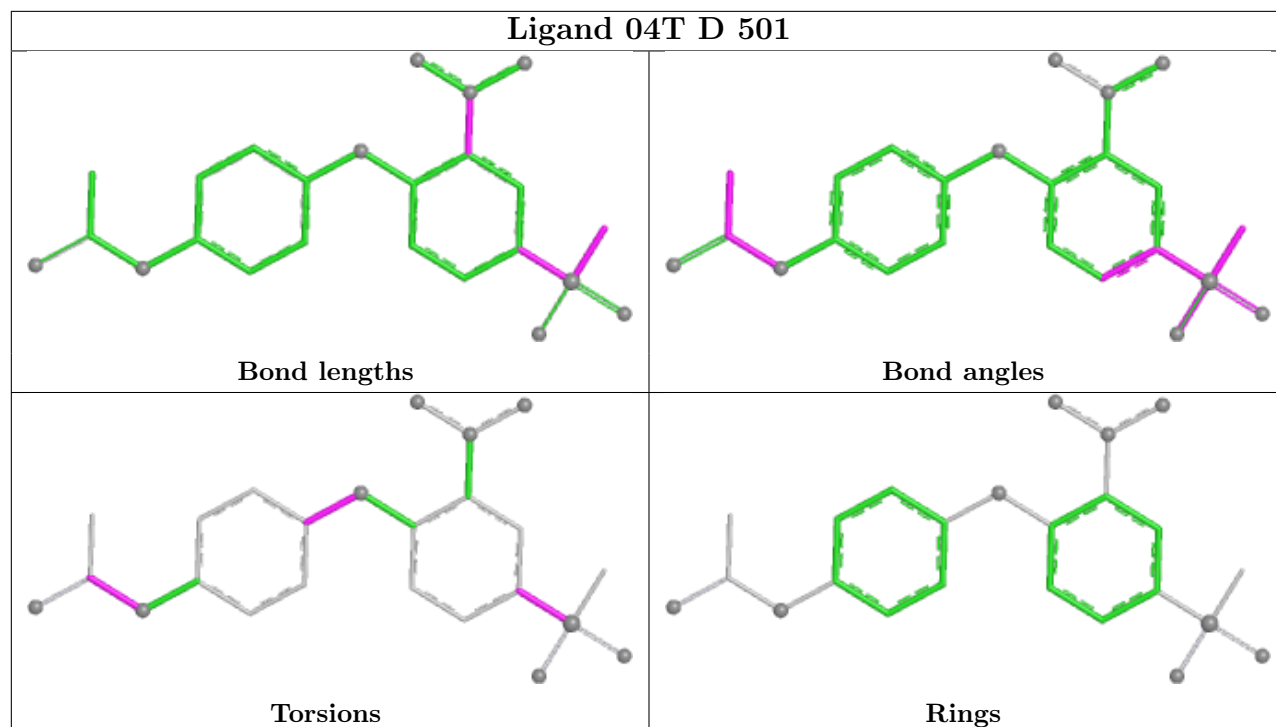
8 monomers are involved in 61 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	501	04T	6	0
2	D	501	04T	6	0
2	F	501	04T	8	0
2	A	501	04T	7	0
2	B	501	04T	9	0
2	E	501	04T	6	0
2	H	501	04T	10	0
2	C	501	04T	9	0

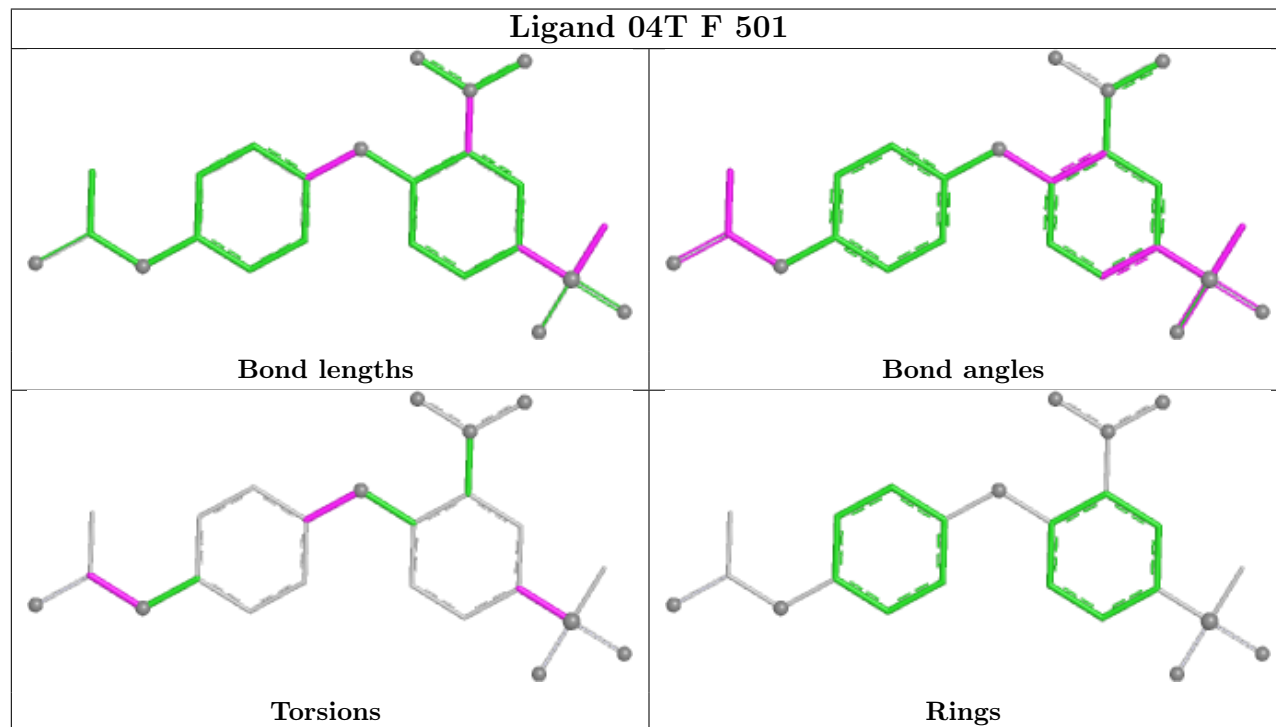
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



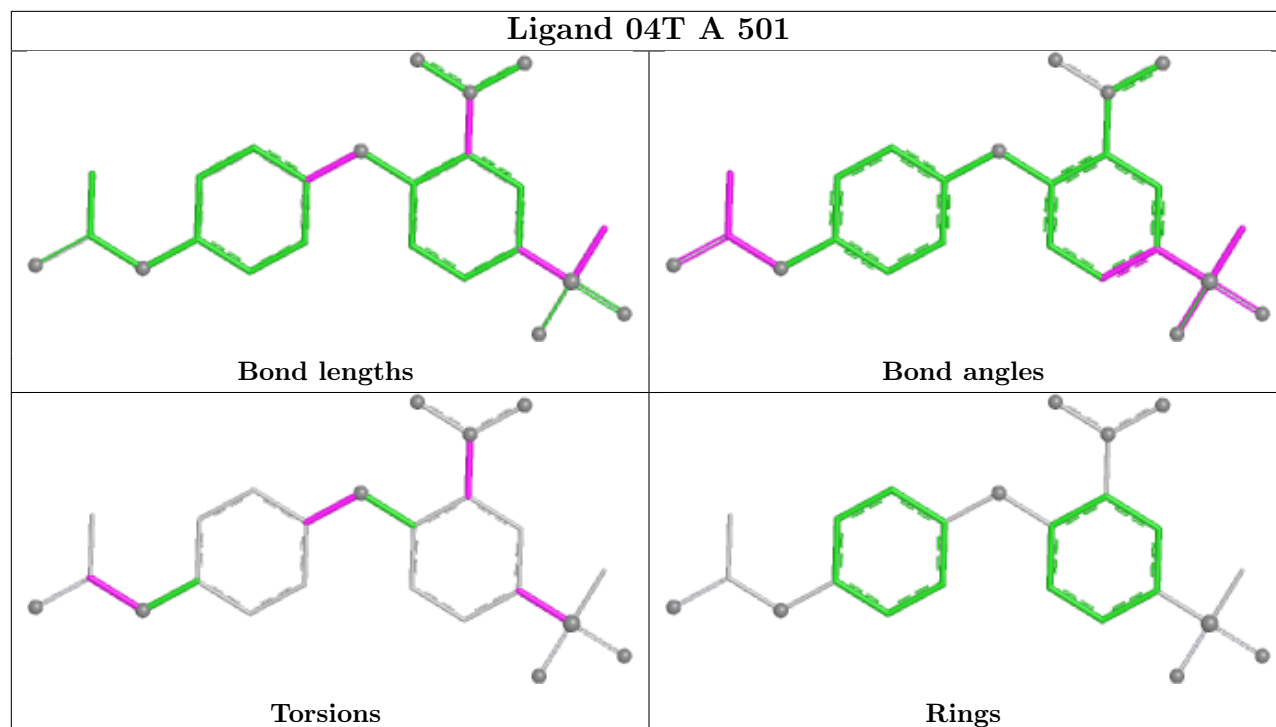
Ligand 04T D 501



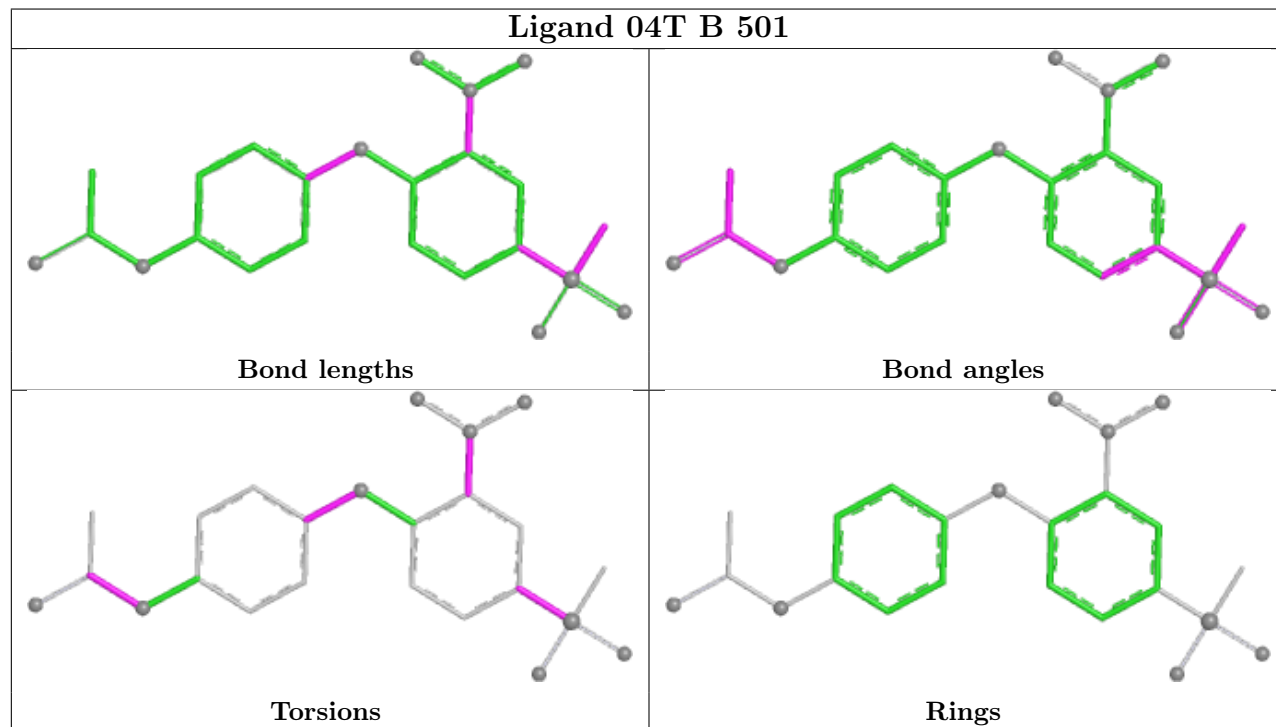
Ligand 04T F 501



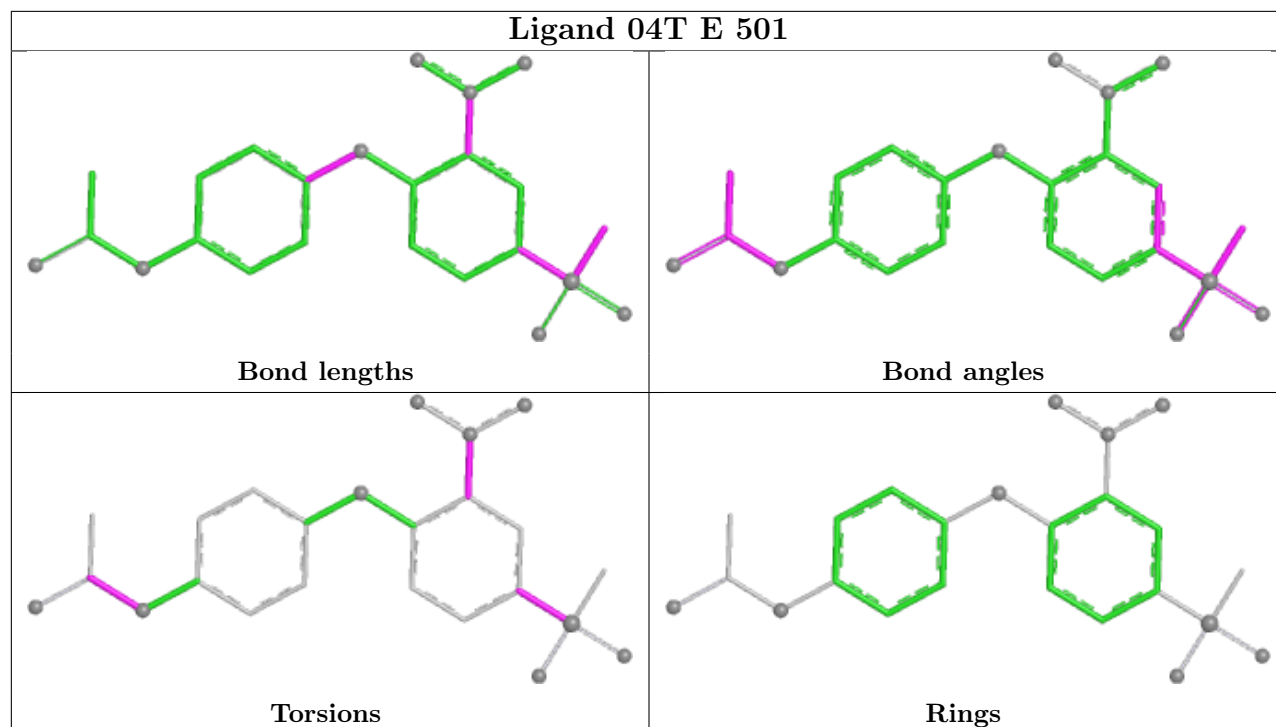
Ligand 04T A 501



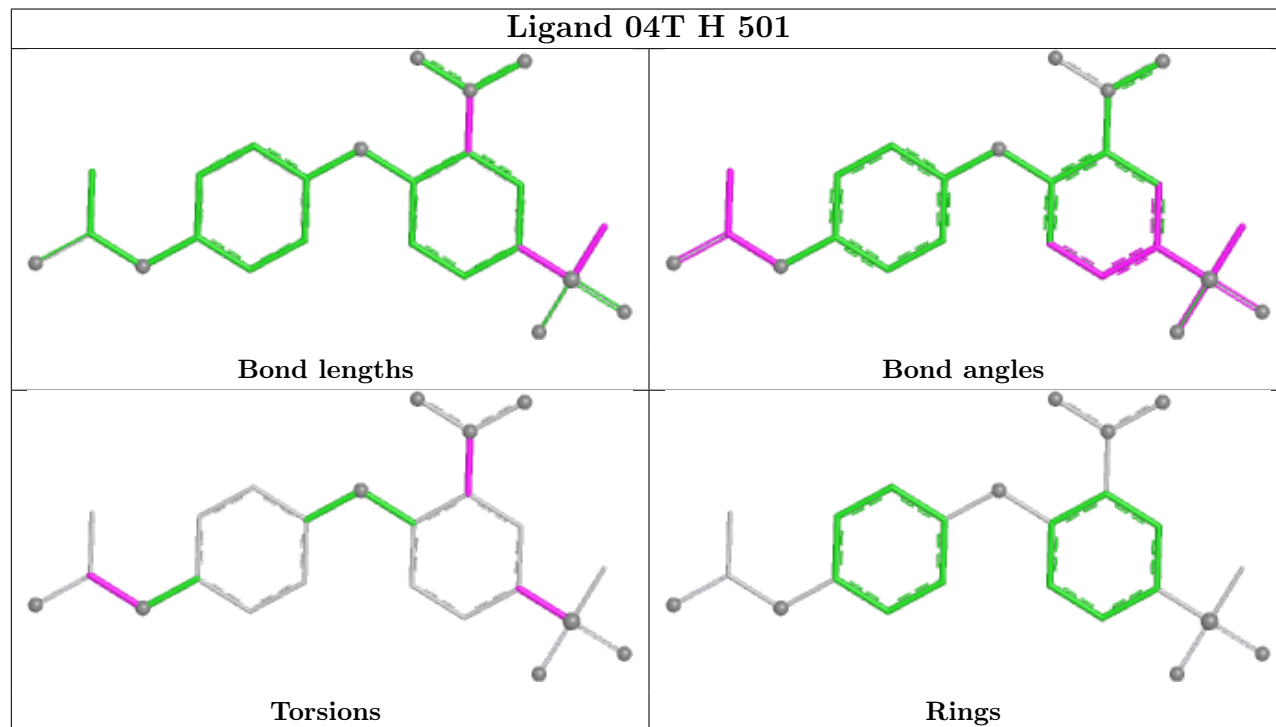
Ligand 04T B 501

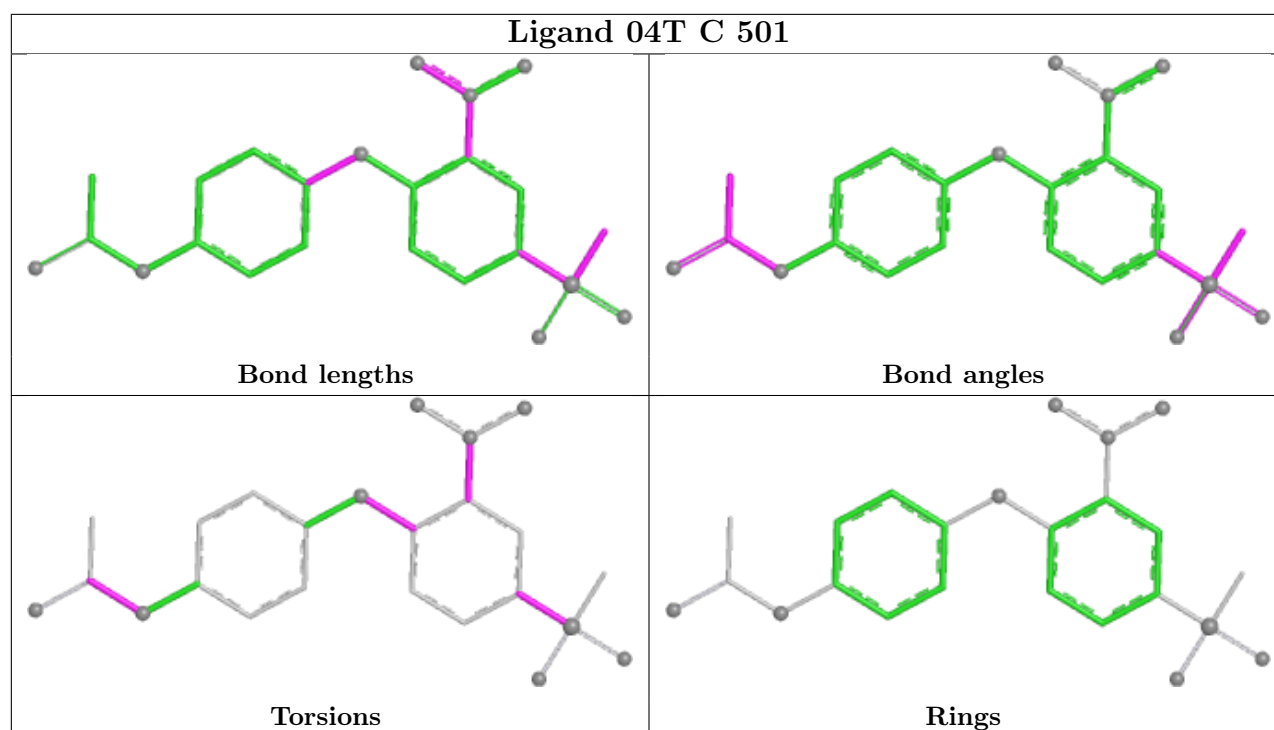


Ligand 04T E 501



Ligand 04T H 501





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	446/469 (95%)	0.42	15 (3%) 48 43	26, 56, 90, 118	3 (0%)
1	B	446/469 (95%)	0.44	9 (2%) 65 61	21, 66, 98, 115	3 (0%)
1	C	447/469 (95%)	0.38	13 (2%) 53 49	23, 64, 115, 132	1 (0%)
1	D	446/469 (95%)	0.14	3 (0%) 84 81	27, 50, 81, 104	2 (0%)
1	E	446/469 (95%)	0.35	5 (1%) 78 75	23, 59, 89, 98	3 (0%)
1	F	446/469 (95%)	0.37	7 (1%) 70 67	28, 69, 118, 134	0
1	G	446/469 (95%)	0.32	10 (2%) 62 58	17, 66, 95, 107	2 (0%)
1	H	447/469 (95%)	0.27	3 (0%) 84 81	28, 58, 95, 128	4 (0%)
All	All	3570/3752 (95%)	0.34	65 (1%) 67 64	17, 60, 102, 134	18 (0%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	93	GLN	3.7
1	C	228	CYS	3.6
1	A	429	VAL	3.4
1	C	447	ALA	3.4
1	F	91	GLN	3.3
1	F	293	VAL	3.1
1	H	447	ALA	3.1
1	G	89	THR	3.1
1	A	291	GLN	3.1
1	G	23	GLN	3.0
1	F	344	VAL	3.0
1	C	227	ALA	2.9
1	A	218	VAL	2.8
1	B	345[A]	ARG	2.8
1	A	430	ARG	2.7
1	B	90	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	G	226	VAL	2.7
1	E	321	THR	2.7
1	A	432	LEU	2.6
1	E	270	PHE	2.6
1	A	433	MET	2.6
1	A	97	LEU	2.6
1	A	310	ALA	2.5
1	B	41	GLN	2.5
1	C	312	THR	2.5
1	C	89	THR	2.5
1	A	224	LEU	2.4
1	C	140	GLU	2.4
1	B	259	ILE	2.4
1	E	322	ASP	2.4
1	G	432	LEU	2.3
1	D	370	ASP	2.3
1	C	294	MET	2.3
1	B	287	ALA	2.3
1	A	446	PRO	2.3
1	E	212	GLY	2.3
1	A	431	PRO	2.2
1	C	287	ALA	2.2
1	F	46	ALA	2.2
1	A	428	LEU	2.2
1	A	225	ASP	2.2
1	A	324	ASP	2.2
1	B	40	GLU	2.2
1	C	218	VAL	2.2
1	G	304	TYR	2.2
1	C	251	CYS	2.2
1	B	140	GLU	2.1
1	F	52	HIS	2.1
1	F	48	ALA	2.1
1	D	93	GLN	2.1
1	G	370	ASP	2.1
1	H	75	GLN	2.1
1	G	49	ALA	2.1
1	B	56	TRP	2.1
1	D	41[A]	GLN	2.0
1	G	91	GLN	2.0
1	B	304	TYR	2.0
1	F	219	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	G	431	PRO	2.0
1	E	433	MET	2.0
1	A	326	GLN	2.0
1	C	338	VAL	2.0
1	H	6	ALA	2.0
1	C	92	THR	2.0
1	C	319	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

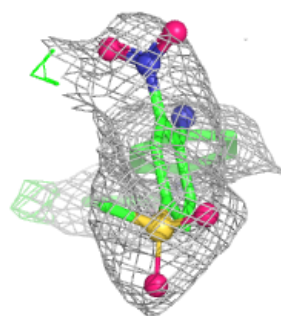
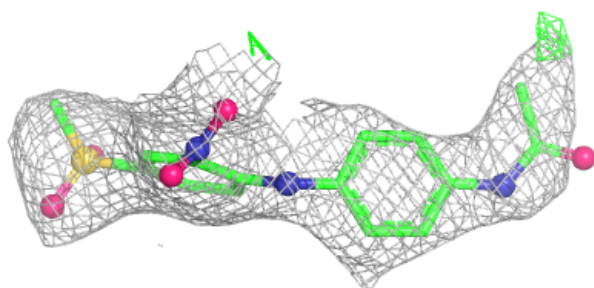
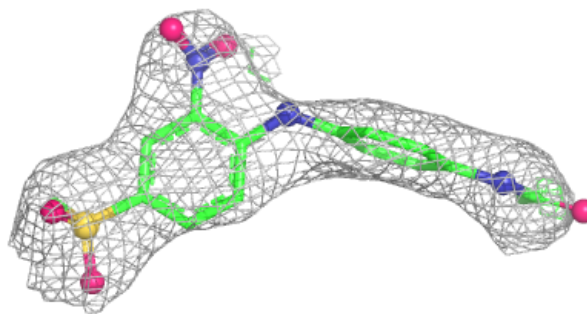
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	04T	B	501	24/24	0.82	0.11	76,78,80,81	0
2	04T	E	501	24/24	0.85	0.12	73,78,80,81	0
2	04T	H	501	24/24	0.86	0.12	62,64,67,68	0
2	04T	A	501	24/24	0.87	0.12	60,66,69,69	0
2	04T	F	501	24/24	0.88	0.11	82,87,90,91	0
2	04T	G	501	24/24	0.88	0.13	76,77,80,81	0
2	04T	D	501	24/24	0.88	0.11	64,66,67,67	0
2	04T	C	501	24/24	0.91	0.11	73,75,77,77	0

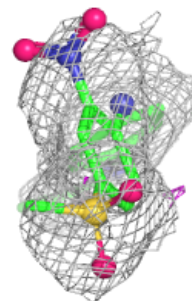
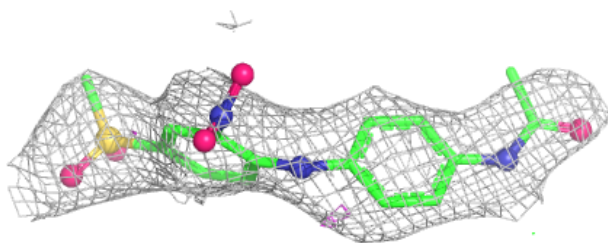
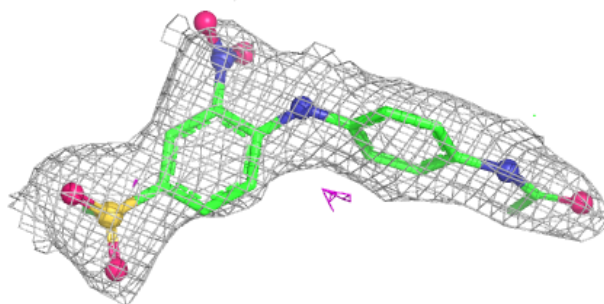
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 04T B 501:

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

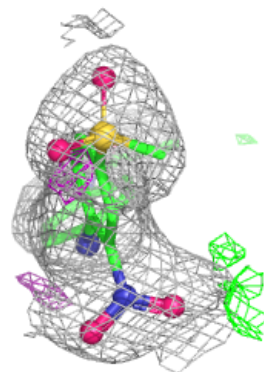
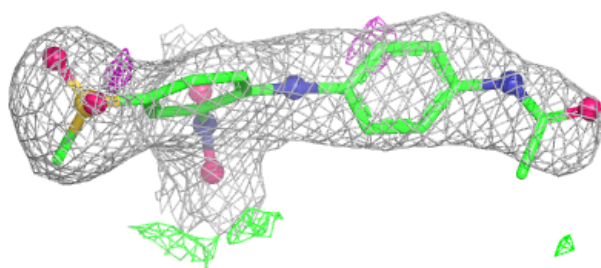
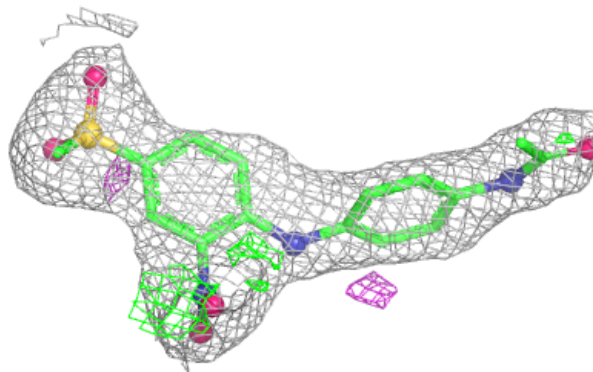
**Electron density around 04T E 501:**

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

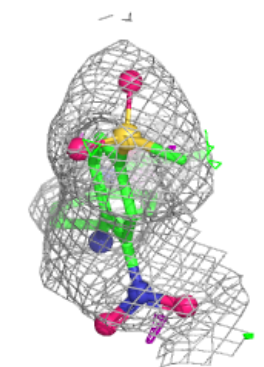
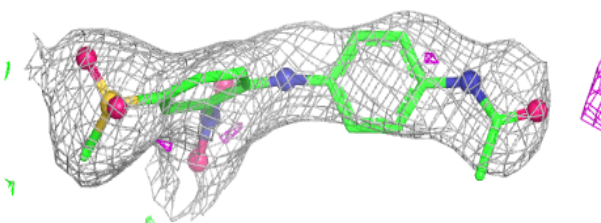
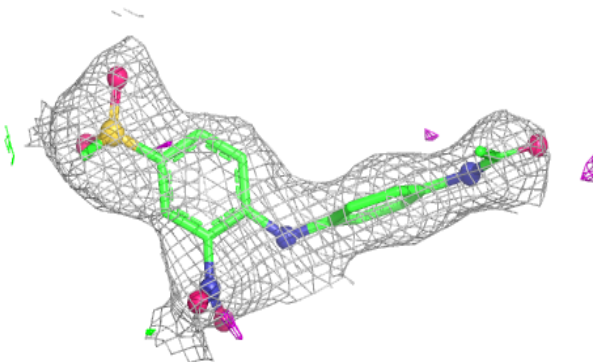


Electron density around 04T H 501:

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

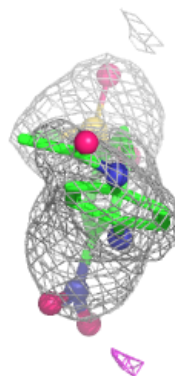
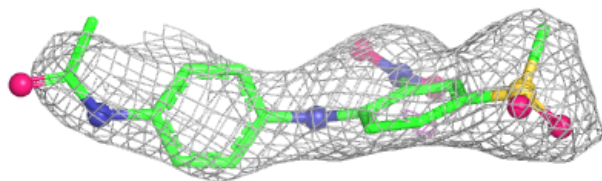
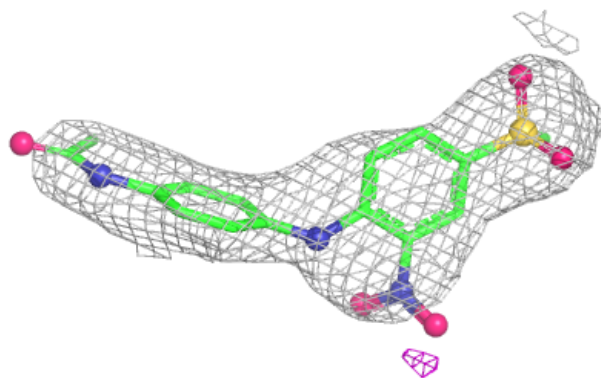
**Electron density around 04T A 501:**

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

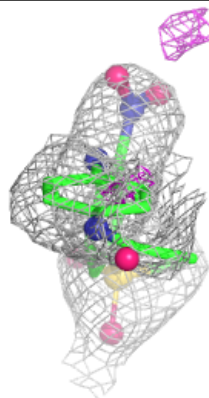
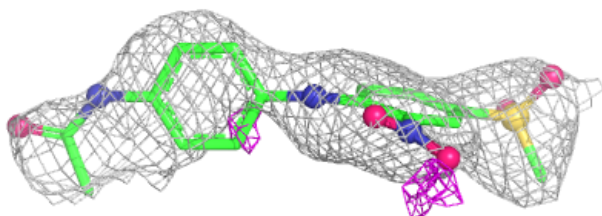
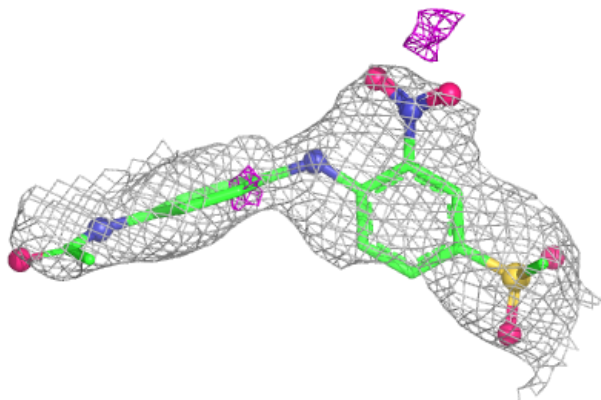


Electron density around 04T F 501:

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

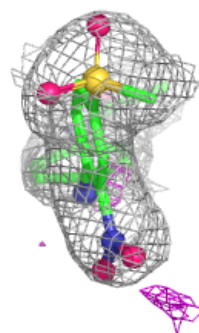
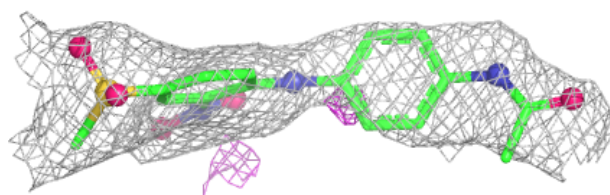
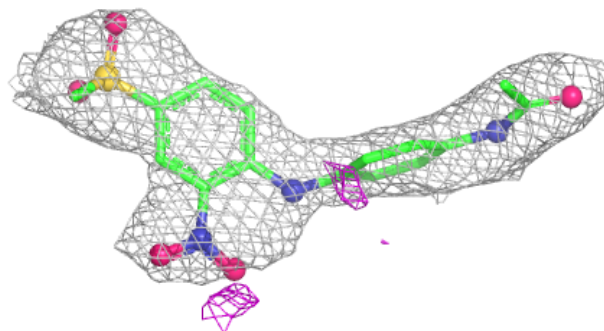
**Electron density around 04T G 501:**

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

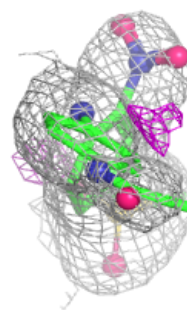
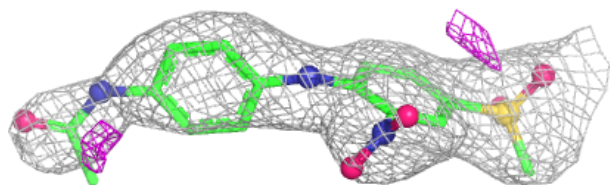
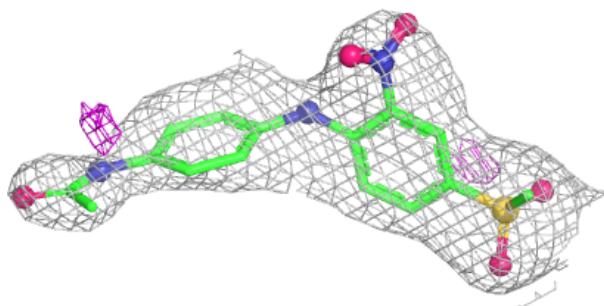


Electron density around 04T D 501:

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

**Electron density around 04T C 501:**

$2mF_o-DF_c$ (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.