Full wwPDB X-ray Structure Validation Report (i)

Mar 10, 2024 – 10:02 AM EDT

PDB ID	:	4HMK
Title	:	Crystal structure of LeuT-E290S with bound Br
Authors	:	Kantcheva, A.K.; Quick, M.; Shi, L.; Winther, A.M.L.; Stolzenberg, S.; Weinstein, H.; Javitch, J.A.; Nissen, P.
Deposited on	:	2012-10-18
Resolution	:	3.00 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:

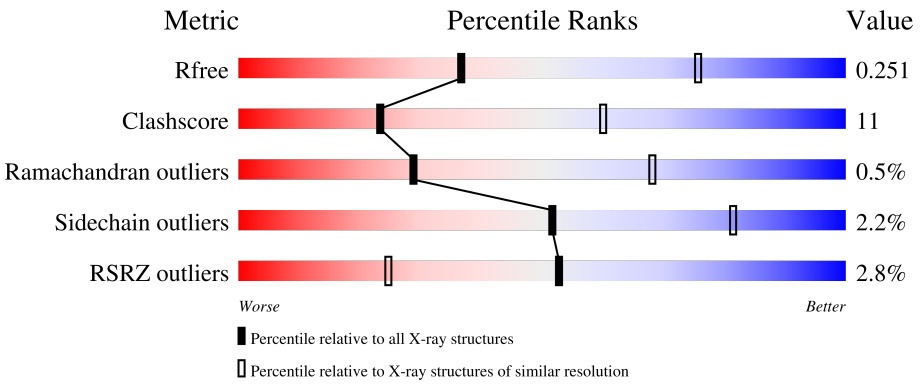
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	513	3% 75% 21% ••
1	B	513	2% 74% 23% ••

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
4	BR	A	611	-	-	X	-
4	BR	B	606	-	-	X	-
4	BR	B	611	-	-	X	-
5	BOG	A	608	-	-	-	X
5	BOG	A	609	-	-	X	X
5	BOG	A	610	-	-	X	X
5	BOG	B	602	-	-	-	X
5	BOG	B	609	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 8240 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.

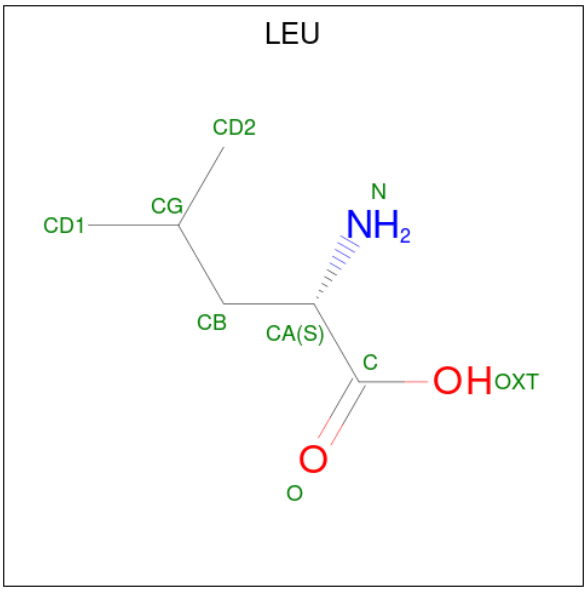
- Molecule 1 is a protein called Transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	501	3987	2700	626	650	11	0	0	0
1	B	506	4027	2728	632	655	12	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	290	SER	GLU	engineered mutation	UNP O67854
B	290	SER	GLU	engineered mutation	UNP O67854

- Molecule 2 is LEUCINE (three-letter code: LEU) (formula: $C_6H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	9	6	1	2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	9	6	1	2	0	0

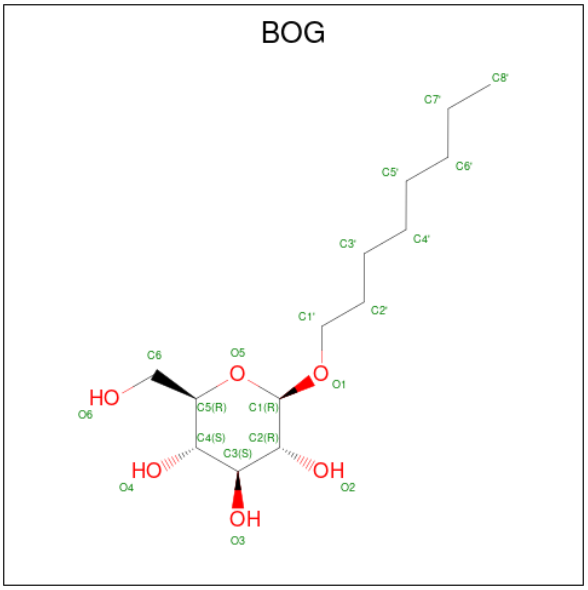
- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
3	A	2	2	2	0	0
			Total	Na		
3	B	2	2	2	0	0

- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Br		
4	A	6	6	6	0	0
			Total	Br		
4	B	5	5	5	0	0

- Molecule 5 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	20	14	6	0	0
			Total	C	O		
5	A	1	20	14	6	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	20	14	6	0	0
			Total	C	O		
5	A	1	20	14	6	0	0
			Total	C	O		
5	A	1	20	14	6	0	0
			Total	C	O		
5	B	1	20	14	6	0	0
			Total	C	O		
5	B	1	20	14	6	0	0
			Total	C	O		
5	B	1	20	14	6	0	0
			Total	C	O		
5	B	1	20	14	6	0	0

- Molecule 6 is water.

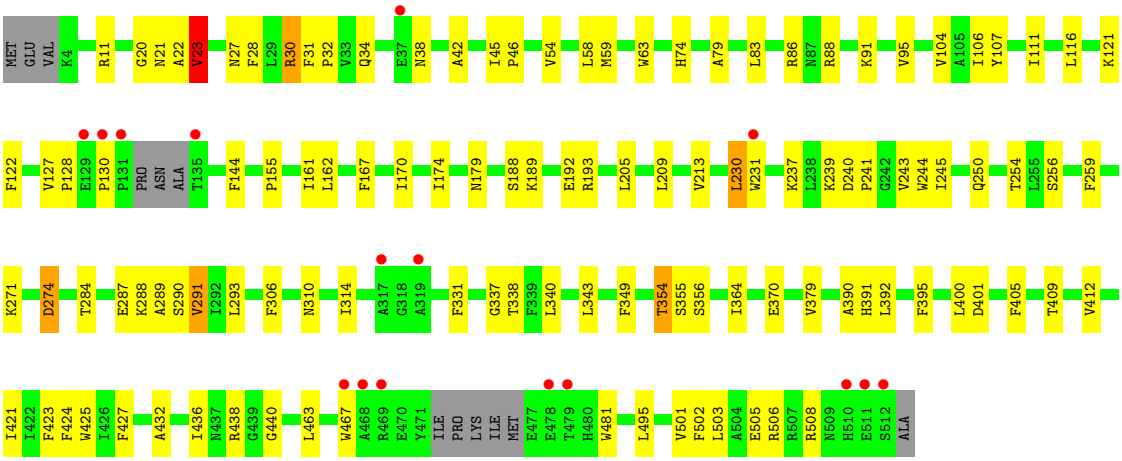
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	7	7	7	0	0
			Total	O		
6	B	6	6	6	0	0

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.

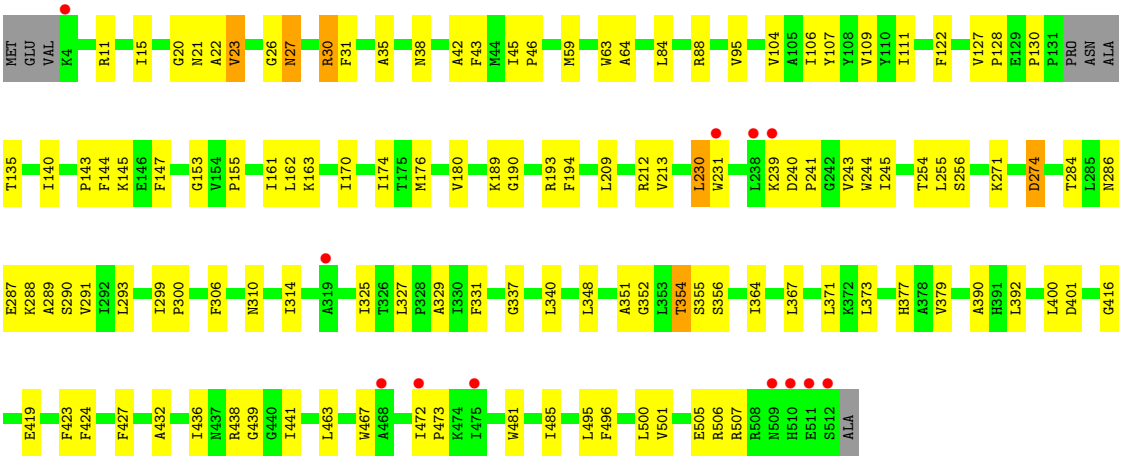
- Molecule 1: Transporter

Chain A:



- Molecule 1: Transporter

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	84.14Å 92.32Å 87.21Å 90.00° 94.25° 90.00°	Depositor
Resolution (Å)	63.30 – 3.00 63.30 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (63.30-3.00) 99.7 (63.30-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.61 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.205 , 0.261 0.201 , 0.251	Depositor DCC
R_{free} test set	1339 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	63.2	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 52.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	8240	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.27% 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: BOG, NA, BR

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	0.42	1/4111 (0.0%)	0.56	1/5596 (0.0%)
1	B	0.40	0/4153	0.55	1/5654 (0.0%)
All	All	0.41	1/8264 (0.0%)	0.55	2/11250 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	23	VAL	CB-CG2	-5.10	1.42	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	351	ALA	C-N-CA	-5.57	110.60	122.30
1	A	354	THR	OG1-CB-CG2	-5.51	97.33	110.00

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	3987	0	4071	95	0
1	B	4027	0	4123	92	0
2	A	9	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	9	0	10	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	6	0	0	4	0
4	B	5	0	0	7	0
5	A	100	0	139	28	0
5	B	80	0	111	10	0
6	A	7	0	0	0	0
6	B	6	0	0	1	0
All	All	8240	0	8464	189	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (189) 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)
5:A:609:BOG:H5	1:B:507:ARG:HA	1.47	0.96
1:A:86:ARG:HH12	5:A:610:BOG:H5'2	1.30	0.93
5:A:608:BOG:H8'3	5:B:601:BOG:H8'2	1.53	0.91
1:A:27:ASN:HD22	1:A:290:SER:HB3	1.38	0.88
1:A:11:ARG:HH22	1:A:274:ASP:HB3	1.41	0.86
5:A:608:BOG:H6'2	5:B:601:BOG:H6'1	1.62	0.81
1:B:11:ARG:HH22	1:B:274:ASP:HB3	1.43	0.81
1:B:364:ILE:HD13	1:B:379:VAL:HG22	1.64	0.78
1:A:86:ARG:NH1	5:A:610:BOG:H5'2	1.97	0.78
1:A:161:ILE:HD12	5:B:601:BOG:H5'1	1.68	0.75
5:A:609:BOG:H5'2	1:B:371:LEU:O	1.88	0.74
1:A:364:ILE:HD13	1:A:379:VAL:HG22	1.70	0.73
1:A:392:LEU:HD13	5:A:608:BOG:H8'1	1.72	0.71
1:B:155:PRO:HD3	1:B:162:LEU:HD23	1.74	0.70
1:A:271:LYS:O	1:A:438:ARG:NH1	2.24	0.70
1:B:135:THR:N	6:B:703:HOH:O	2.26	0.69
1:B:27:ASN:HD22	1:B:290:SER:HB3	1.58	0.69
1:A:189:LYS:HA	1:A:193:ARG:HG3	1.77	0.67
1:A:83:LEU:HD21	5:A:610:BOG:H4'2	1.76	0.67
1:B:256:SER:HG	2:B:603:LEU:N	1.95	0.65
5:A:609:BOG:H8'1	4:B:611:BR:BR	2.53	0.64
1:A:155:PRO:HD3	1:A:162:LEU:HD23	1.79	0.64
1:A:245:ILE:HG12	1:A:463:LEU:HD12	1.79	0.64
1:A:256:SER:HG	2:A:601:LEU:N	1.96	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:367:LEU:HB3	1:B:373:LEU:HD22	1.81	0.63
1:A:11:ARG:HH22	1:A:274:ASP:CB	2.11	0.61
1:B:45:ILE:HB	1:B:46:PRO:HD3	1.82	0.61
1:A:88:ARG:NH1	1:A:91:LYS:HE3	2.15	0.61
1:B:11:ARG:HH22	1:B:274:ASP:CB	2.15	0.60
1:B:11:ARG:NH2	1:B:274:ASP:HB3	2.15	0.59
1:B:189:LYS:HA	1:B:193:ARG:HG3	1.83	0.59
1:B:392:LEU:HD13	5:B:609:BOG:H8'1	1.84	0.58
1:B:241:PRO:O	1:B:245:ILE:HG13	2.03	0.58
1:A:503:LEU:HD22	5:A:609:BOG:H4'2	1.85	0.58
1:A:338:THR:HB	4:A:613:BR:BR	2.58	0.58
1:A:23:VAL:HG12	1:A:23:VAL:O	2.04	0.57
1:A:111:ILE:HD11	1:A:400:LEU:HD11	1.86	0.57
1:A:45:ILE:HB	1:A:46:PRO:HD3	1.86	0.57
1:B:31:PHE:CD1	4:B:606:BR:BR	3.13	0.57
1:A:11:ARG:NH2	1:A:274:ASP:HB3	2.15	0.56
1:B:245:ILE:HG12	1:B:463:LEU:HD12	1.88	0.56
1:A:331:PHE:O	1:A:337:GLY:HA3	2.06	0.56
1:A:88:ARG:HH11	1:A:88:ARG:HA	1.69	0.56
1:A:31:PHE:HB2	4:A:604:BR:BR	2.61	0.55
1:B:377:HIS:ND1	4:B:611:BR:BR	2.87	0.55
1:B:31:PHE:HB2	4:B:606:BR:BR	2.62	0.54
1:A:395:PHE:HB3	5:B:601:BOG:H5'2	1.90	0.54
1:A:34:GLN:NE2	4:A:611:BR:BR	2.95	0.54
1:B:23:VAL:O	1:B:23:VAL:HG12	2.06	0.54
1:A:506:ARG:HG3	5:A:609:BOG:H8'3	1.89	0.54
1:B:212:ARG:HH11	1:B:212:ARG:HG3	1.73	0.54
1:A:74:HIS:NE2	5:A:610:BOG:H3'1	2.24	0.53
1:A:130:PRO:HG3	1:A:144:PHE:HE2	1.72	0.53
1:B:254:THR:OG1	4:B:606:BR:BR	2.79	0.53
1:A:27:ASN:ND2	1:A:290:SER:HB3	2.18	0.52
1:A:241:PRO:O	1:A:245:ILE:HG13	2.09	0.52
1:B:289:ALA:O	1:B:293:LEU:HB2	2.10	0.52
1:A:27:ASN:HB3	1:A:290:SER:HB3	1.91	0.51
5:A:609:BOG:H8'2	1:B:373:LEU:HD12	1.93	0.51
1:A:167:PHE:HE1	5:A:606:BOG:H6'2	1.76	0.50
1:B:255:LEU:HD23	1:B:286:ASN:HD22	1.75	0.50
1:A:30:ARG:HD3	4:A:611:BR:BR	2.67	0.50
1:A:239:LYS:HD2	1:A:239:LYS:N	2.25	0.50
1:B:95:VAL:HG22	1:B:501:VAL:HG21	1.93	0.50
1:B:271:LYS:O	1:B:438:ARG:NH1	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:PRO:HB3	1:A:244:TRP:CE3	2.47	0.50
1:A:370:GLU:OE1	5:A:610:BOG:H61	2.11	0.50
1:A:88:ARG:HD3	1:A:91:LYS:HD2	1.94	0.49
1:A:167:PHE:HD1	5:A:606:BOG:H4'2	1.77	0.49
1:B:239:LYS:N	1:B:239:LYS:HD2	2.27	0.49
1:B:88:ARG:HH11	1:B:88:ARG:HA	1.77	0.49
1:B:230:LEU:HB3	1:B:231:TRP:CD1	2.47	0.49
1:B:284:THR:HG22	1:B:288:LYS:HE3	1.94	0.49
1:A:423:PHE:O	1:A:427:PHE:HB3	2.12	0.49
1:A:506:ARG:CB	5:A:609:BOG:H6'2	2.43	0.49
1:A:495:LEU:CB	5:B:602:BOG:H61	2.42	0.49
1:B:432:ALA:O	1:B:436:ILE:HG13	2.13	0.48
1:A:495:LEU:HB2	5:B:602:BOG:H61	1.96	0.48
1:B:240:ASP:HB3	1:B:243:VAL:HG23	1.96	0.48
1:B:27:ASN:HB3	1:B:290:SER:HB3	1.96	0.48
1:B:287:GLU:O	1:B:291:VAL:HG13	2.13	0.48
1:B:416:GLY:O	1:B:419:GLU:HG2	2.14	0.48
1:B:127:VAL:HB	1:B:128:PRO:HD2	1.96	0.48
1:A:79:ALA:HA	5:A:610:BOG:H2	1.96	0.47
1:B:46:PRO:HB3	1:B:244:TRP:CE3	2.49	0.47
1:B:104:VAL:HA	1:B:107:TYR:CE2	2.49	0.47
1:B:505:GLU:HG3	1:B:506:ARG:NH1	2.30	0.47
1:B:495:LEU:HB3	5:B:602:BOG:H8'2	1.95	0.47
1:A:179:ASN:OD1	1:A:356:SER:OG	2.30	0.47
1:B:30:ARG:HD3	4:B:610:BR:BR	2.70	0.47
1:A:130:PRO:HG3	1:A:144:PHE:CE2	2.49	0.47
1:B:245:ILE:HG12	1:B:463:LEU:CD1	2.45	0.47
1:B:163:LYS:NZ	5:B:609:BOG:H61	2.30	0.47
1:A:505:GLU:HG3	1:A:506:ARG:NH1	2.30	0.47
1:A:240:ASP:HB3	1:A:243:VAL:HG23	1.96	0.46
1:A:250:GLN:O	1:A:254:THR:HG22	2.15	0.46
1:B:231:TRP:HH2	1:B:291:VAL:HB	1.80	0.46
1:B:11:ARG:O	1:B:15:ILE:HG12	2.16	0.46
1:A:506:ARG:HB2	5:A:609:BOG:H6'2	1.97	0.46
1:A:122:PHE:CD1	1:A:127:VAL:HG11	2.50	0.46
1:B:122:PHE:CD1	1:B:127:VAL:HG11	2.51	0.46
5:A:609:BOG:C8'	4:B:611:BR:BR	3.18	0.46
1:A:306:PHE:HB3	1:A:310:ASN:HB2	1.98	0.46
1:B:111:ILE:HD11	1:B:400:LEU:HD11	1.98	0.46
1:A:245:ILE:HG12	1:A:463:LEU:CD1	2.46	0.45
1:A:20:GLY:O	1:A:21:ASN:C	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:MET:O	1:A:63:TRP:HD1	1.99	0.45
1:A:287:GLU:O	1:A:291:VAL:HG13	2.17	0.45
1:B:331:PHE:O	1:B:337:GLY:HA3	2.16	0.45
1:A:205:LEU:HD22	1:A:343:LEU:HD22	1.98	0.45
1:A:432:ALA:O	1:A:436:ILE:HG13	2.16	0.45
1:A:83:LEU:CD2	5:A:610:BOG:H4'2	2.44	0.45
5:A:609:BOG:H5	1:B:507:ARG:CA	2.34	0.45
1:B:153:GLY:HA3	1:B:163:LYS:O	2.17	0.45
1:B:255:LEU:HD23	1:B:286:ASN:ND2	2.31	0.45
1:B:306:PHE:HB3	1:B:310:ASN:HB2	1.98	0.45
1:B:472:ILE:N	1:B:473:PRO:HD2	2.32	0.45
1:B:140:ILE:O	1:B:143:PRO:HD2	2.17	0.45
1:A:128:PRO:HG2	1:A:144:PHE:CE1	2.50	0.45
1:B:255:LEU:CD2	1:B:286:ASN:HD22	2.29	0.45
1:B:109:VAL:HG22	1:B:356:SER:HB2	1.99	0.44
1:A:161:ILE:HD11	1:B:481:TRP:CE3	2.51	0.44
1:A:106:ILE:HG23	1:A:390:ALA:HA	1.99	0.44
1:B:310:ASN:O	1:B:314:ILE:HG13	2.17	0.44
1:A:340:LEU:HD12	1:A:340:LEU:HA	1.75	0.44
1:A:230:LEU:HB3	1:A:231:TRP:CD1	2.53	0.44
1:A:259:PHE:CZ	1:A:412:VAL:HG11	2.53	0.44
1:A:502:PHE:O	1:A:506:ARG:HG2	2.17	0.44
1:A:167:PHE:CE1	5:A:606:BOG:H6'2	2.53	0.44
1:A:237:LYS:HB3	1:A:243:VAL:HG21	2.00	0.44
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.53	0.44
1:B:38:ASN:O	1:B:42:ALA:HB3	2.17	0.44
1:B:423:PHE:O	1:B:427:PHE:HB3	2.17	0.44
1:B:35:ALA:HA	1:B:43:PHE:CD2	2.53	0.43
1:B:46:PRO:HG2	1:B:243:VAL:HG12	1.99	0.43
1:B:170:ILE:O	1:B:174:ILE:HG13	2.18	0.43
1:B:144:PHE:CE2	1:B:329:ALA:HA	2.53	0.43
1:A:274:ASP:OD1	1:A:440:GLY:HA2	2.19	0.43
1:A:370:GLU:HB3	5:A:610:BOG:H62	2.01	0.43
1:B:26:GLY:HA3	2:B:603:LEU:OXT	2.18	0.43
1:B:145:LYS:HA	1:B:325:ILE:HD13	2.01	0.43
1:B:106:ILE:HG23	1:B:390:ALA:HA	2.00	0.43
1:A:161:ILE:HD11	1:B:481:TRP:CZ3	2.54	0.43
1:A:505:GLU:HB2	1:A:508:ARG:HH21	1.84	0.43
1:A:245:ILE:HG23	1:A:463:LEU:HD11	2.00	0.43
1:B:153:GLY:HA3	1:B:163:LYS:C	2.39	0.43
1:B:340:LEU:HD12	1:B:340:LEU:HA	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ALA:O	1:A:293:LEU:HB2	2.19	0.42
1:B:348:LEU:O	1:B:352:GLY:N	2.52	0.42
1:B:194:PHE:HE2	1:B:354:THR:HG23	1.85	0.42
1:B:128:PRO:HD3	1:B:147:PHE:CD1	2.54	0.42
1:A:391:HIS:HD2	5:A:607:BOG:H3'1	1.85	0.42
1:A:481:TRP:CE3	1:B:161:ILE:HD11	2.55	0.42
1:B:11:ARG:HH21	1:B:441:ILE:HD12	1.85	0.42
1:A:481:TRP:CZ3	1:B:161:ILE:HD11	2.55	0.42
1:B:84:LEU:HB3	1:B:427:PHE:HB2	2.02	0.42
1:B:496:PHE:CZ	1:B:500:LEU:HD11	2.55	0.42
1:A:188:SER:O	1:A:192:GLU:HB3	2.20	0.42
5:A:609:BOG:H1'1	1:B:507:ARG:HB2	2.01	0.41
1:B:299:ILE:HB	1:B:300:PRO:HD3	2.02	0.41
5:A:607:BOG:H3'2	5:A:607:BOG:H6'1	1.90	0.41
1:B:194:PHE:CE2	1:B:354:THR:HG23	2.55	0.41
1:A:54:VAL:O	1:A:58:LEU:HB2	2.20	0.41
1:A:405:PHE:O	1:A:409:THR:HB	2.19	0.41
1:A:116:LEU:HB2	1:A:349:PHE:CD1	2.56	0.41
1:A:284:THR:HG22	1:A:288:LYS:HE3	2.02	0.41
5:A:610:BOG:H3'2	5:A:610:BOG:H7'2	2.01	0.41
1:A:310:ASN:O	1:A:314:ILE:HG13	2.19	0.41
1:B:176:MET:O	1:B:180:VAL:HG23	2.20	0.41
1:B:485:ILE:HG12	5:B:601:BOG:H8'3	2.03	0.41
1:A:46:PRO:HG2	1:A:243:VAL:HG12	2.02	0.41
1:A:421:ILE:O	1:A:425:TRP:HB2	2.21	0.41
1:B:64:ALA:HB2	1:B:436:ILE:HG12	2.03	0.41
1:B:209:LEU:O	1:B:213:VAL:HG23	2.21	0.41
1:A:95:VAL:HG22	1:A:501:VAL:HG21	2.03	0.41
1:B:274:ASP:HB2	1:B:439:GLY:O	2.21	0.41
1:B:327:LEU:HA	1:B:327:LEU:HD12	1.83	0.41
1:B:212:ARG:HG3	1:B:212:ARG:NH1	2.36	0.41
1:B:20:GLY:O	1:B:21:ASN:C	2.55	0.40
1:A:28:PHE:O	1:A:32:PRO:HG2	2.21	0.40
1:A:161:ILE:HD13	1:A:161:ILE:HA	1.87	0.40
1:A:38:ASN:O	1:A:42:ALA:HB3	2.21	0.40
1:A:170:ILE:O	1:A:174:ILE:HG13	2.21	0.40
1:B:59:MET:O	1:B:63:TRP:HD1	2.04	0.40
1:A:23:VAL:O	1:A:23:VAL:CG1	2.67	0.40
1:A:205:LEU:CD2	1:A:343:LEU:HD22	2.51	0.40
1:A:209:LEU:O	1:A:213:VAL:HG23	2.21	0.40
1:B:130:PRO:HG3	1:B:144:PHE:HE2	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:LYS:HD2	1:A:167:PHE:CE2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	495/513 (96%)	475 (96%)	18 (4%)	2 (0%)	34	72
1	B	502/513 (98%)	469 (93%)	30 (6%)	3 (1%)	25	64
All	All	997/1026 (97%)	944 (95%)	48 (5%)	5 (0%)	29	68

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	23	VAL
1	A	23	VAL
1	B	22	ALA
1	A	22	ALA
1	B	190	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	405/415 (98%)	396 (98%)	9 (2%)	52 81
1	B	410/415 (99%)	401 (98%)	9 (2%)	52 81
All	All	815/830 (98%)	797 (98%)	18 (2%)	52 81

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ARG
1	A	230	LEU
1	A	274	ASP
1	A	291	VAL
1	A	354	THR
1	A	355	SER
1	A	401	ASP
1	A	424	PHE
1	A	467	TRP
1	B	27	ASN
1	B	30	ARG
1	B	230	LEU
1	B	274	ASP
1	B	354	THR
1	B	355	SER
1	B	401	ASP
1	B	424	PHE
1	B	467	TRP

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 26 ligands modelled in this entry, 15 are monoatomic - leaving 11 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	$\#\|Z\| > 2$
5	BOG	A	607	-	20,20,20	1.32	3 (15%)	25,25,25	1.49	4 (16%)
2	LEU	A	601	3	7,8,8	1.27	1 (14%)	9,10,10	1.68	2 (22%)
5	BOG	A	606	-	20,20,20	1.35	3 (15%)	25,25,25	1.21	3 (12%)
5	BOG	A	610	-	20,20,20	2.41	7 (35%)	25,25,25	2.11	6 (24%)
2	LEU	B	603	3	7,8,8	1.08	1 (14%)	9,10,10	1.67	2 (22%)
5	BOG	B	609	-	20,20,20	1.39	3 (15%)	25,25,25	1.10	3 (12%)
5	BOG	B	608	-	20,20,20	1.41	3 (15%)	25,25,25	1.45	3 (12%)
5	BOG	B	601	-	20,20,20	1.36	3 (15%)	25,25,25	1.36	4 (16%)
5	BOG	A	609	-	20,20,20	1.46	3 (15%)	25,25,25	1.29	4 (16%)
5	BOG	B	602	-	20,20,20	1.77	5 (25%)	25,25,25	2.78	9 (36%)
5	BOG	A	608	-	20,20,20	1.32	3 (15%)	25,25,25	1.21	4 (16%)

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
5	BOG	A	607	-	-	6/11/31/31	0/1/1/1
2	LEU	A	601	3	-	0/8/8/8	-
5	BOG	A	606	-	-	2/11/31/31	0/1/1/1
5	BOG	A	610	-	-	8/11/31/31	0/1/1/1
2	LEU	B	603	3	-	2/8/8/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BOG	B	609	-	-	4/11/31/31	0/1/1/1
5	BOG	B	608	-	-	3/11/31/31	0/1/1/1
5	BOG	B	601	-	-	5/11/31/31	0/1/1/1
5	BOG	A	609	-	-	7/11/31/31	0/1/1/1
5	BOG	B	602	-	-	10/11/31/31	0/1/1/1
5	BOG	A	608	-	-	4/11/31/31	0/1/1/1

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	610	BOG	O5-C1	5.08	1.54	1.41
5	A	610	BOG	O1-C1	-5.03	1.31	1.40
5	A	610	BOG	O2-C2	-4.71	1.31	1.43
5	B	602	BOG	O5-C1	4.54	1.53	1.41
5	B	601	BOG	O5-C1	4.14	1.52	1.41
5	B	608	BOG	O5-C1	4.06	1.52	1.41
5	A	609	BOG	O5-C1	3.95	1.51	1.41
5	B	609	BOG	O5-C1	3.95	1.51	1.41
5	A	610	BOG	C3-C2	-3.93	1.42	1.52
5	A	606	BOG	O5-C1	3.90	1.51	1.41
5	A	607	BOG	O5-C1	3.89	1.51	1.41
5	A	608	BOG	O5-C1	3.58	1.51	1.41
5	B	602	BOG	C3-C2	-3.24	1.44	1.52
5	A	609	BOG	O1-C1	-3.20	1.34	1.40
2	A	601	LEU	OXT-C	-3.14	1.20	1.30
5	A	608	BOG	O1-C1	-2.95	1.35	1.40
5	B	609	BOG	O1-C1	-2.93	1.35	1.40
5	B	602	BOG	O3-C3	-2.90	1.36	1.43
5	A	610	BOG	C4-C3	-2.77	1.45	1.52
5	A	606	BOG	O1-C1	-2.74	1.35	1.40
2	B	603	LEU	OXT-C	-2.74	1.21	1.30
5	A	610	BOG	C1-C2	-2.64	1.44	1.52
5	A	608	BOG	C3-C2	-2.60	1.45	1.52
5	B	608	BOG	O1-C1	-2.55	1.35	1.40
5	A	609	BOG	C3-C2	-2.52	1.45	1.52
5	B	608	BOG	C3-C2	-2.50	1.46	1.52
5	B	609	BOG	C3-C2	-2.48	1.46	1.52
5	B	602	BOG	C4-C5	-2.45	1.47	1.53
5	A	607	BOG	O1-C1	-2.44	1.36	1.40
5	A	606	BOG	C3-C2	-2.43	1.46	1.52
5	B	601	BOG	O1-C1	-2.41	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	602	BOG	C4-C3	-2.37	1.46	1.52
5	A	610	BOG	O5-C5	-2.35	1.38	1.44
5	A	607	BOG	C3-C2	-2.26	1.46	1.52
5	B	601	BOG	C3-C2	-2.24	1.46	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	602	BOG	C1'-O1-C1	7.79	126.76	113.84
5	A	610	BOG	O2-C2-C1	-6.73	93.69	110.05
5	A	607	BOG	C1'-O1-C1	5.24	122.52	113.84
5	B	608	BOG	C1'-O1-C1	5.18	122.42	113.84
5	B	602	BOG	O3-C3-C4	-5.16	98.41	110.35
5	B	602	BOG	O3-C3-C2	-5.12	98.52	110.35
5	B	602	BOG	C6-C5-C4	-4.87	101.60	113.00
2	B	603	LEU	OXT-C-O	-4.06	114.87	124.09
2	A	601	LEU	OXT-C-O	-4.04	114.91	124.09
5	B	602	BOG	O5-C5-C6	3.74	115.74	106.44
5	A	610	BOG	C3-C4-C5	-3.73	103.59	110.24
5	A	610	BOG	O2-C2-C3	-3.57	102.08	110.35
5	B	601	BOG	C1'-O1-C1	3.29	119.29	113.84
5	B	609	BOG	C1'-O1-C1	3.24	119.21	113.84
5	A	609	BOG	O5-C5-C6	3.16	114.29	106.44
5	A	610	BOG	C1-C2-C3	3.06	116.36	110.00
5	A	608	BOG	C1'-O1-C1	3.01	118.84	113.84
5	A	609	BOG	C1'-O1-C1	2.99	118.80	113.84
5	A	608	BOG	O1-C1-C2	2.85	112.76	108.30
5	B	602	BOG	O6-C6-C5	2.65	120.37	111.29
5	A	609	BOG	C1-O5-C5	-2.63	108.52	113.69
5	B	601	BOG	O1-C1'-C2'	2.59	118.66	109.56
5	A	610	BOG	O4-C4-C5	2.58	115.69	109.30
5	B	602	BOG	O1-C1'-C2'	2.57	118.58	109.56
5	A	606	BOG	O1-C1-C2	2.56	112.30	108.30
5	A	610	BOG	O1-C1'-C2'	2.41	117.99	109.56
5	A	606	BOG	O1-C1'-C2'	2.37	117.86	109.56
5	B	602	BOG	O2-C2-C1	2.35	115.75	110.05
5	A	609	BOG	O5-C5-C4	2.33	113.93	109.69
5	A	606	BOG	C1'-O1-C1	2.32	117.69	113.84
2	B	603	LEU	OXT-C-CA	2.31	121.26	113.38
5	B	601	BOG	C4-C3-C2	-2.30	106.81	110.82
5	B	608	BOG	C6-C5-C4	-2.25	107.73	113.00
5	A	607	BOG	O1-C1-C2	2.20	111.74	108.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	607	BOG	O1-C1'-C2'	2.19	117.24	109.56
5	A	608	BOG	O6-C6-C5	2.19	118.80	111.29
5	B	602	BOG	O1-C1-C2	2.18	111.70	108.30
2	A	601	LEU	OXT-C-CA	2.14	120.67	113.38
5	A	608	BOG	O1-C1'-C2'	2.12	116.98	109.56
5	B	609	BOG	O1-C1'-C2'	2.10	116.93	109.56
5	A	607	BOG	C1-O5-C5	-2.09	109.58	113.69
5	B	609	BOG	O6-C6-C5	2.07	118.39	111.29
5	B	608	BOG	C4-C3-C2	-2.06	107.22	110.82
5	B	601	BOG	C6-C5-C4	-2.00	108.32	113.00

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	607	BOG	O5-C1-O1-C1'
5	A	608	BOG	O5-C1-O1-C1'
5	A	609	BOG	C2'-C1'-O1-C1
5	A	610	BOG	C2-C1-O1-C1'
5	A	610	BOG	O5-C1-O1-C1'
5	B	601	BOG	C2'-C1'-O1-C1
5	B	602	BOG	C2-C1-O1-C1'
5	B	602	BOG	O5-C1-O1-C1'
5	A	609	BOG	C4-C5-C6-O6
5	A	609	BOG	O5-C5-C6-O6
5	A	609	BOG	O1-C1'-C2'-C3'
5	B	602	BOG	C4-C5-C6-O6
5	B	602	BOG	O5-C5-C6-O6
5	A	610	BOG	O1-C1'-C2'-C3'
5	B	608	BOG	C3'-C4'-C5'-C6'
5	B	608	BOG	C4-C5-C6-O6
5	A	610	BOG	C2'-C3'-C4'-C5'
5	A	609	BOG	C3'-C4'-C5'-C6'
5	B	602	BOG	C4'-C5'-C6'-C7'
5	B	602	BOG	C1'-C2'-C3'-C4'
5	B	601	BOG	C3'-C4'-C5'-C6'
5	A	606	BOG	C3'-C4'-C5'-C6'
5	A	607	BOG	C3'-C4'-C5'-C6'
5	A	609	BOG	C1'-C2'-C3'-C4'
5	B	608	BOG	O5-C5-C6-O6
5	A	608	BOG	C3'-C4'-C5'-C6'
5	B	602	BOG	C2'-C3'-C4'-C5'

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Mol	Chain	Res	Type	Atoms
5	A	610	BOG	C4'-C5'-C6'-C7'
5	A	607	BOG	O5-C5-C6-O6
5	B	602	BOG	C3'-C4'-C5'-C6'
5	A	607	BOG	O1-C1'-C2'-C3'
5	B	601	BOG	O1-C1'-C2'-C3'
5	B	609	BOG	O5-C1-O1-C1'
5	A	607	BOG	C4-C5-C6-O6
5	A	608	BOG	C2-C1-O1-C1'
5	A	610	BOG	C5'-C6'-C7'-C8'
5	A	610	BOG	C2'-C1'-O1-C1
5	B	602	BOG	C2'-C1'-O1-C1
5	B	609	BOG	O1-C1'-C2'-C3'
5	A	606	BOG	O1-C1'-C2'-C3'
5	B	609	BOG	C5'-C6'-C7'-C8'
5	A	610	BOG	C1'-C2'-C3'-C4'
2	B	603	LEU	CA-CB-CG-CD2
5	A	607	BOG	C4'-C5'-C6'-C7'
5	A	609	BOG	C5'-C6'-C7'-C8'
5	B	601	BOG	C2'-C3'-C4'-C5'
5	B	609	BOG	C3'-C4'-C5'-C6'
5	A	608	BOG	C5'-C6'-C7'-C8'
5	B	601	BOG	C1'-C2'-C3'-C4'
5	B	602	BOG	O1-C1'-C2'-C3'
2	B	603	LEU	OXT-C-CA-N

There are no ring outliers.

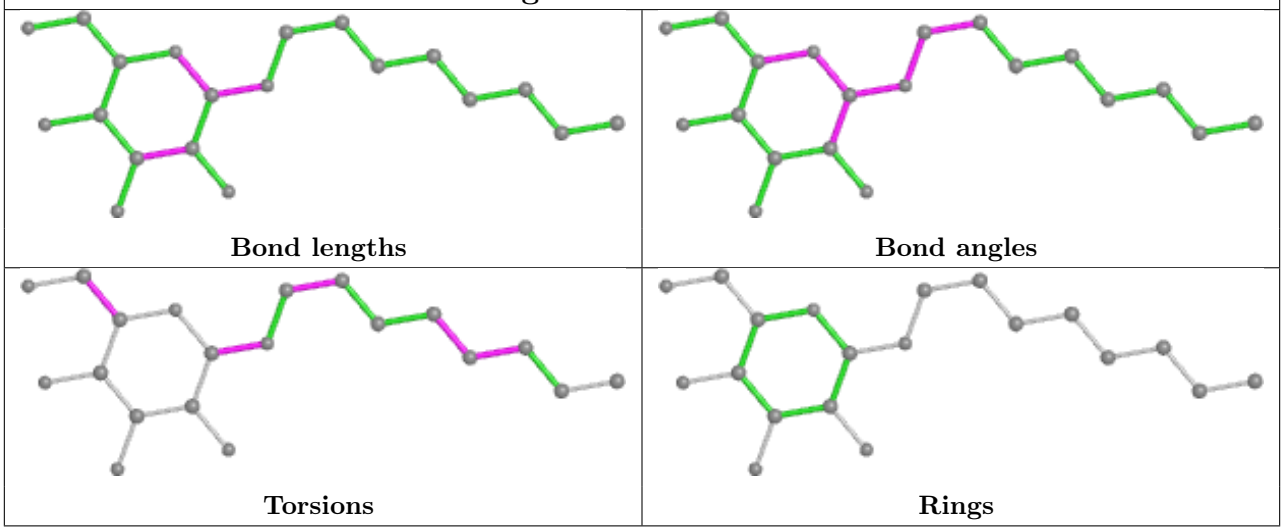
10 monomers are involved in 39 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	607	BOG	2	0
2	A	601	LEU	1	0
5	A	606	BOG	3	0
5	A	610	BOG	9	0
2	B	603	LEU	2	0
5	B	609	BOG	2	0
5	B	601	BOG	5	0
5	A	609	BOG	11	0
5	B	602	BOG	3	0
5	A	608	BOG	3	0

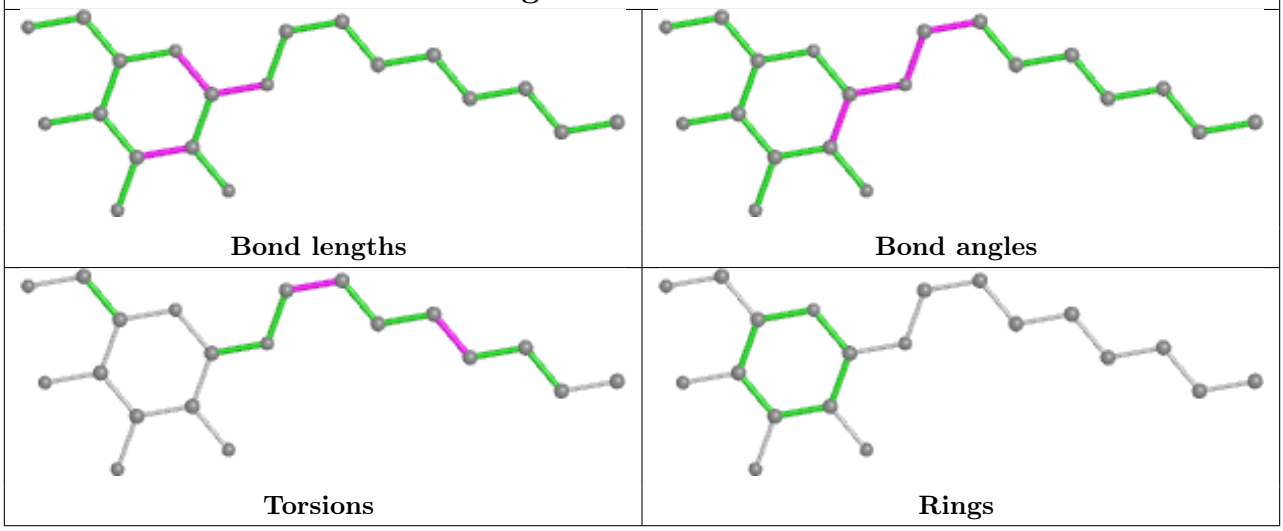
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.

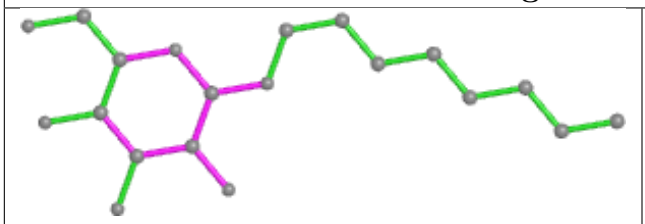
Ligand BOG A 607



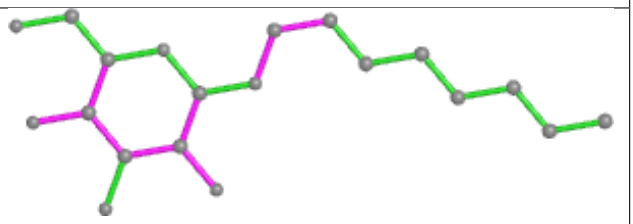
Ligand BOG A 606



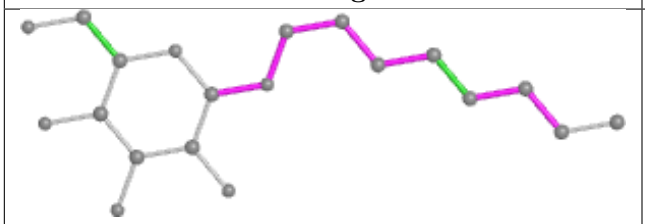
Ligand BOG A 610



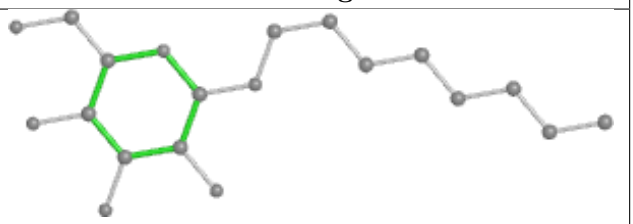
Bond lengths



Bond angles

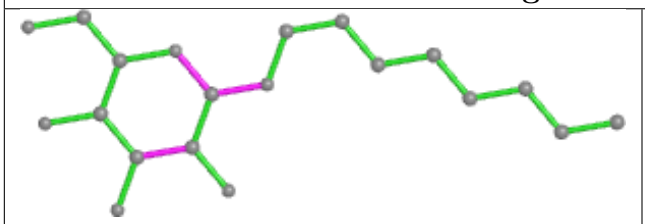


Torsions

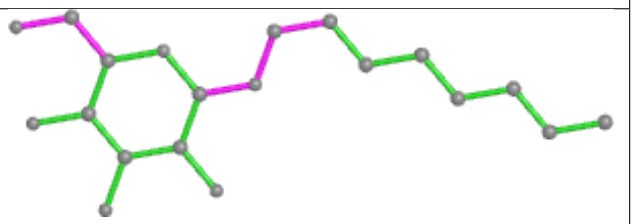


Rings

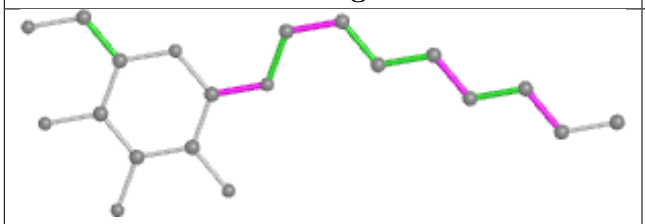
Ligand BOG B 609



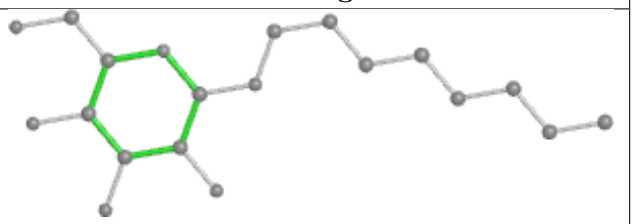
Bond lengths



Bond angles

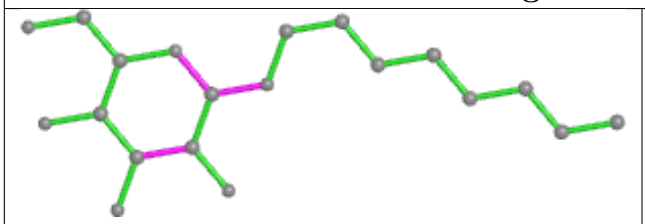


Torsions

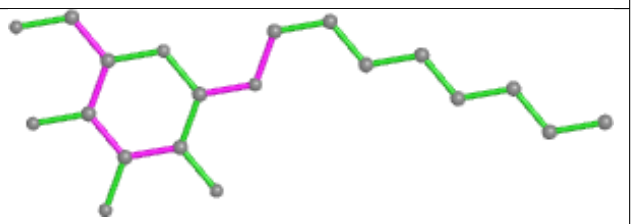


Rings

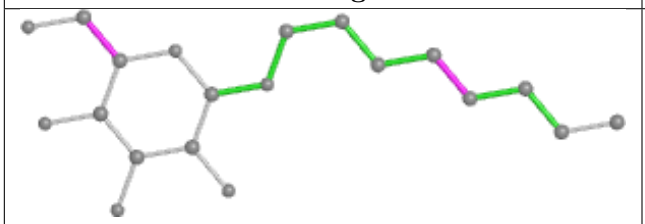
Ligand BOG B 608



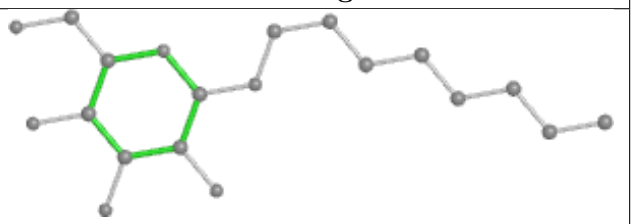
Bond lengths



Bond angles

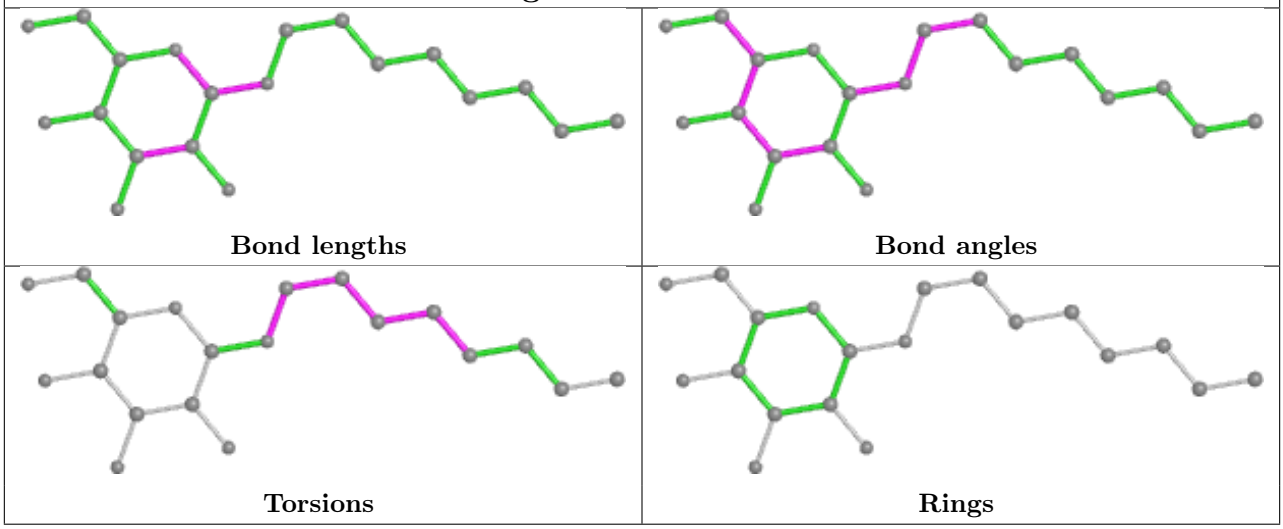


Torsions

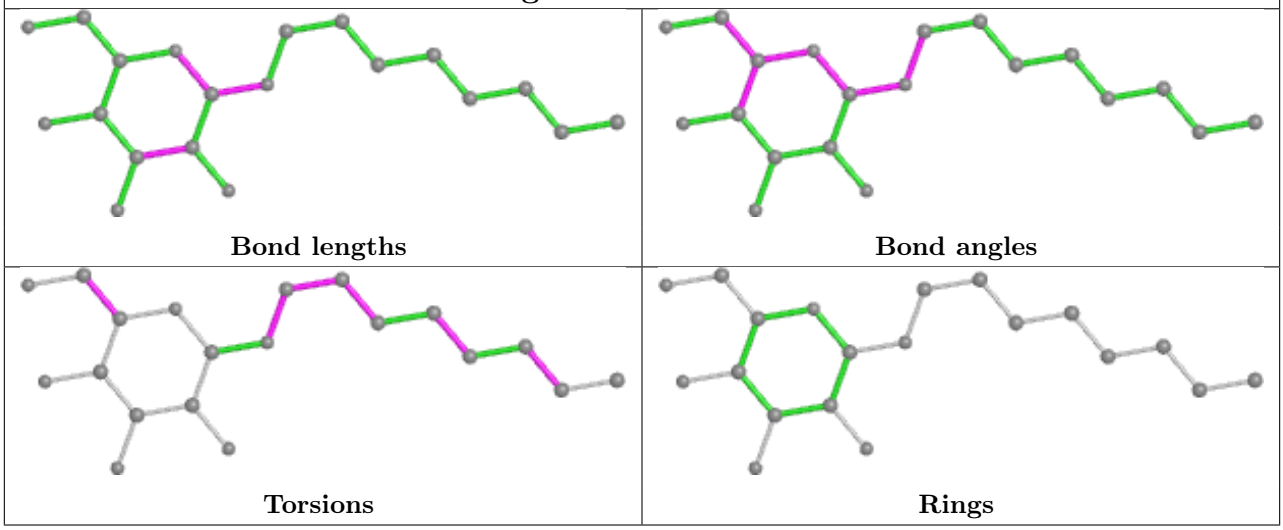


Rings

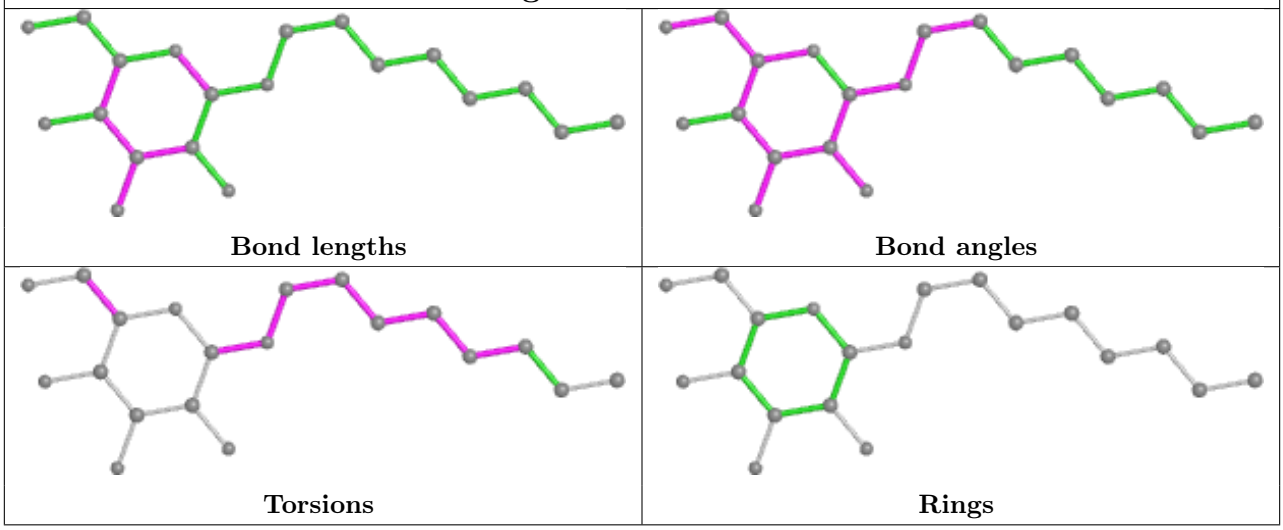
Ligand BOG B 601

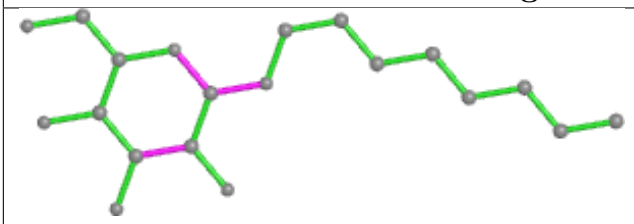
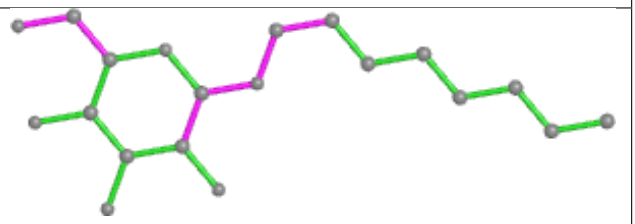
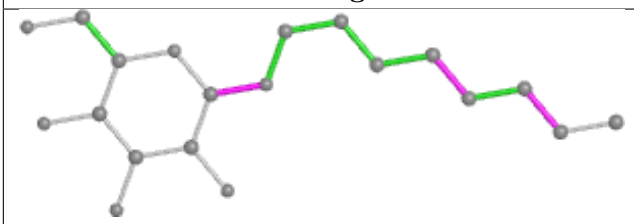
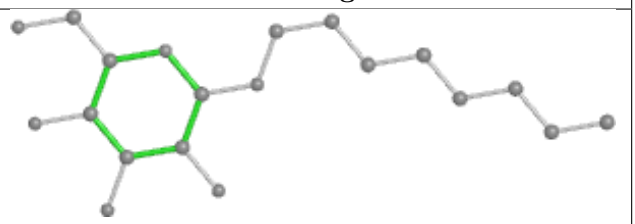


Ligand BOG A 609



Ligand BOG B 602



Ligand BOG A 608	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	501/513 (97%)	-0.08	16 (3%)	47	20	34, 50, 87, 110	0
1	B	506/513 (98%)	-0.13	12 (2%)	59	30	34, 52, 90, 110	0
All	All	1007/1026 (98%)	-0.10	28 (2%)	53	25	34, 51, 88, 110	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	PRO	7.7
1	B	512	SER	6.0
1	B	511	GLU	5.4
1	B	468	ALA	5.1
1	A	512	SER	4.6
1	A	511	GLU	4.2
1	A	317	ALA	4.0
1	B	475	ILE	4.0
1	B	472	ILE	3.6
1	B	239	LYS	3.3
1	A	135	THR	3.1
1	A	319	ALA	3.0
1	A	478	GLU	3.0
1	B	510	HIS	2.8
1	A	469	ARG	2.7
1	A	467	TRP	2.7
1	A	129	GLU	2.6
1	A	468	ALA	2.5
1	A	479	THR	2.4
1	B	319	ALA	2.3
1	A	130	PRO	2.3
1	A	231	TRP	2.2
1	B	4	LYS	2.2
1	B	238	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	510	HIS	2.1
1	B	509	ASN	2.1
1	A	37	GLU	2.1
1	B	231	TRP	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 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
4	BR	A	613	1/1	0.10	0.36	237,237,237,237	0
5	BOG	B	609	20/20	0.62	0.48	34,71,80,81	20
5	BOG	A	609	20/20	0.66	0.69	49,69,85,88	20
5	BOG	B	602	20/20	0.67	0.41	41,54,69,77	20
3	NA	A	603	1/1	0.68	0.25	46,46,46,46	0
5	BOG	A	610	20/20	0.71	0.45	39,84,97,97	20
5	BOG	A	608	20/20	0.73	0.48	48,65,100,104	20
5	BOG	A	606	20/20	0.79	0.33	47,79,95,100	20
3	NA	A	602	1/1	0.80	0.36	40,40,40,40	0
3	NA	B	604	1/1	0.81	0.22	49,49,49,49	0
5	BOG	B	601	20/20	0.81	0.35	44,74,92,94	20
5	BOG	B	608	20/20	0.85	0.36	28,61,78,81	20
5	BOG	A	607	20/20	0.92	0.25	50,68,80,84	0
4	BR	B	612	1/1	0.93	0.09	90,90,90,90	1
4	BR	A	614	1/1	0.93	0.10	63,63,63,63	1
4	BR	B	610	1/1	0.95	0.12	88,88,88,88	1
4	BR	B	611	1/1	0.95	0.08	84,84,84,84	1
4	BR	A	612	1/1	0.95	0.11	96,96,96,96	1
3	NA	B	605	1/1	0.95	0.27	55,55,55,55	0

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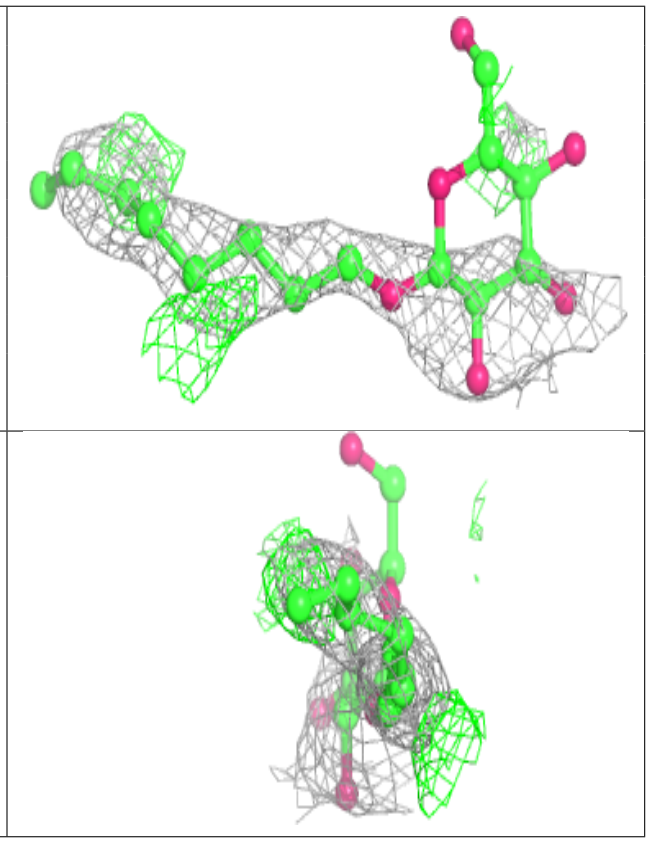
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	BR	A	604	1/1	0.95	0.09	73,73,73,73	1
4	BR	B	607	1/1	0.95	0.07	75,75,75,75	1
4	BR	A	605	1/1	0.98	0.06	59,59,59,59	1
4	BR	B	606	1/1	0.98	0.14	77,77,77,77	1
4	BR	A	611	1/1	0.98	0.16	76,76,76,76	1
2	LEU	B	603	9/9	0.98	0.27	38,45,49,54	0
2	LEU	A	601	9/9	0.98	0.25	33,42,44,45	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.

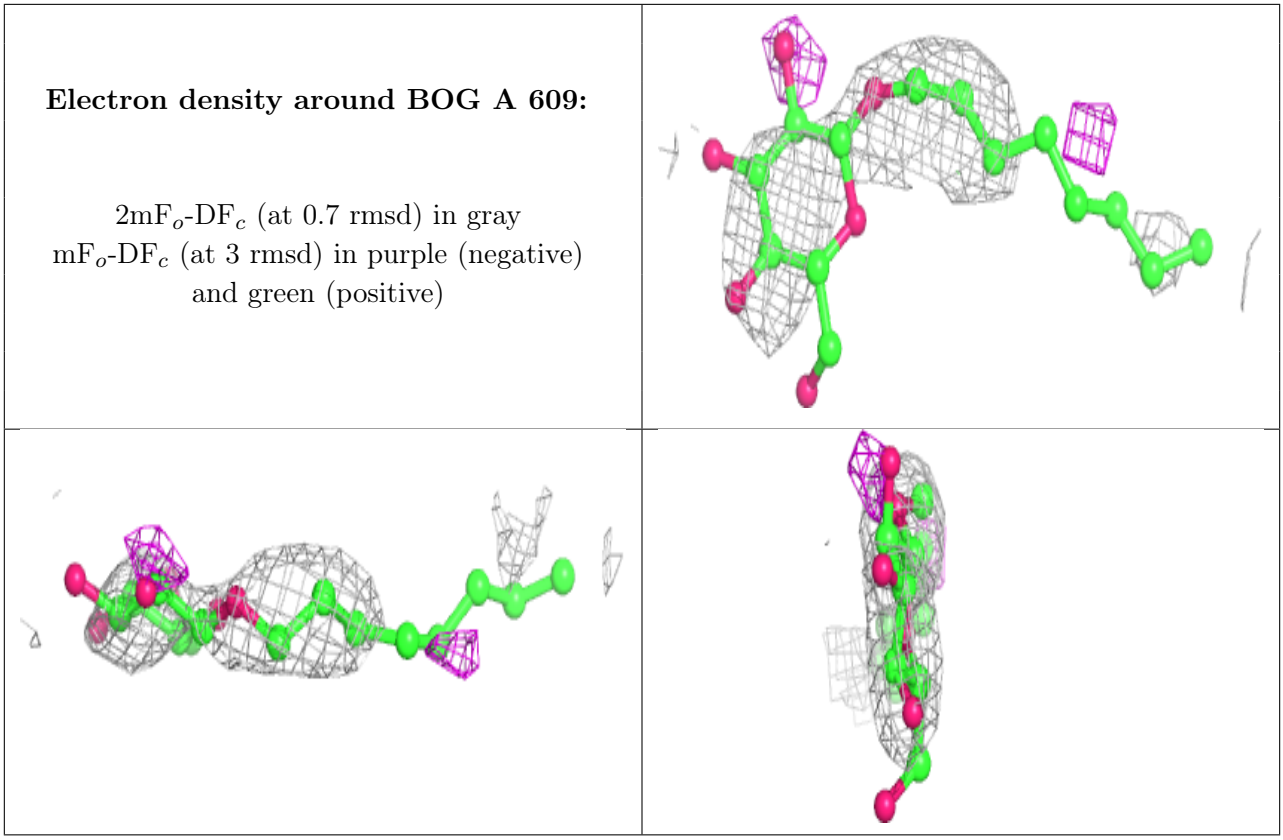
Electron density around BOG B 609:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)



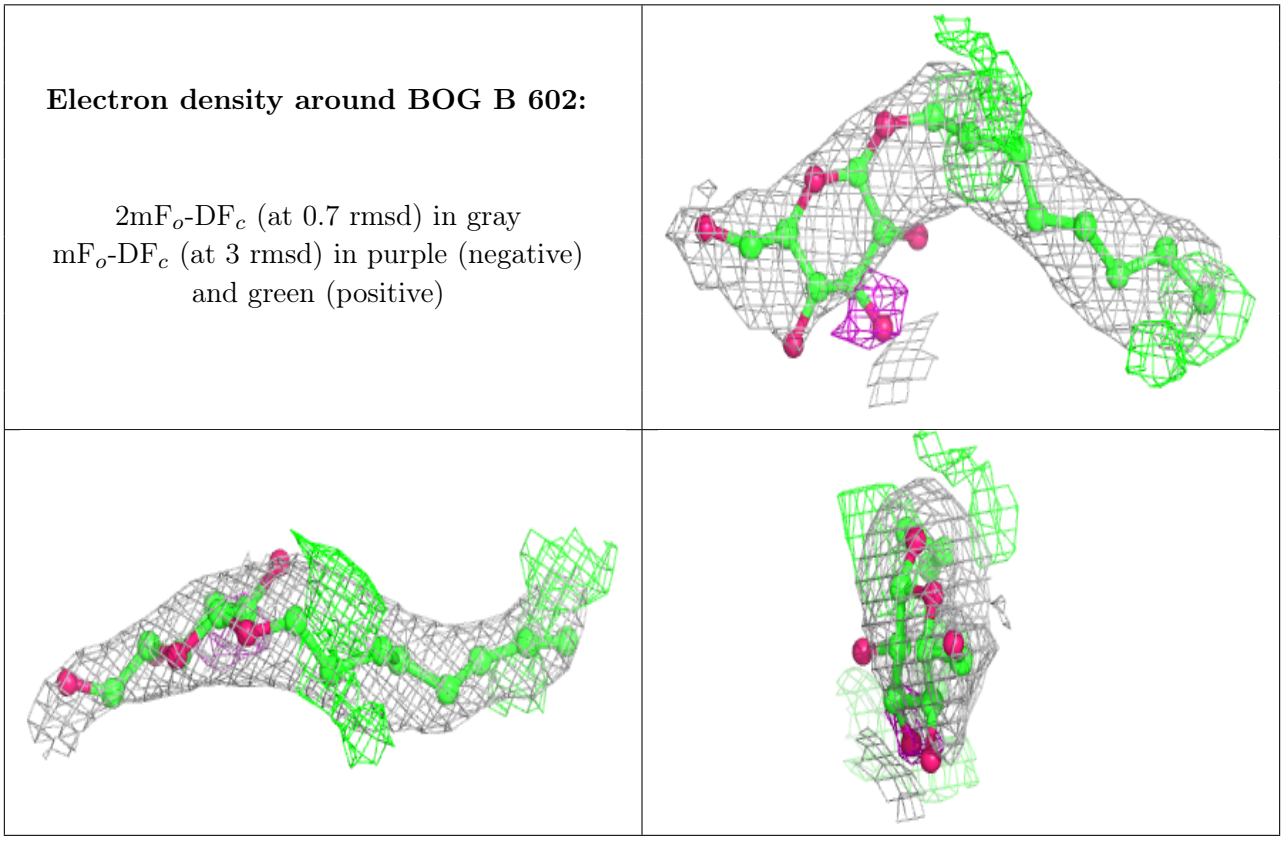
Electron density around BOG A 609:

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



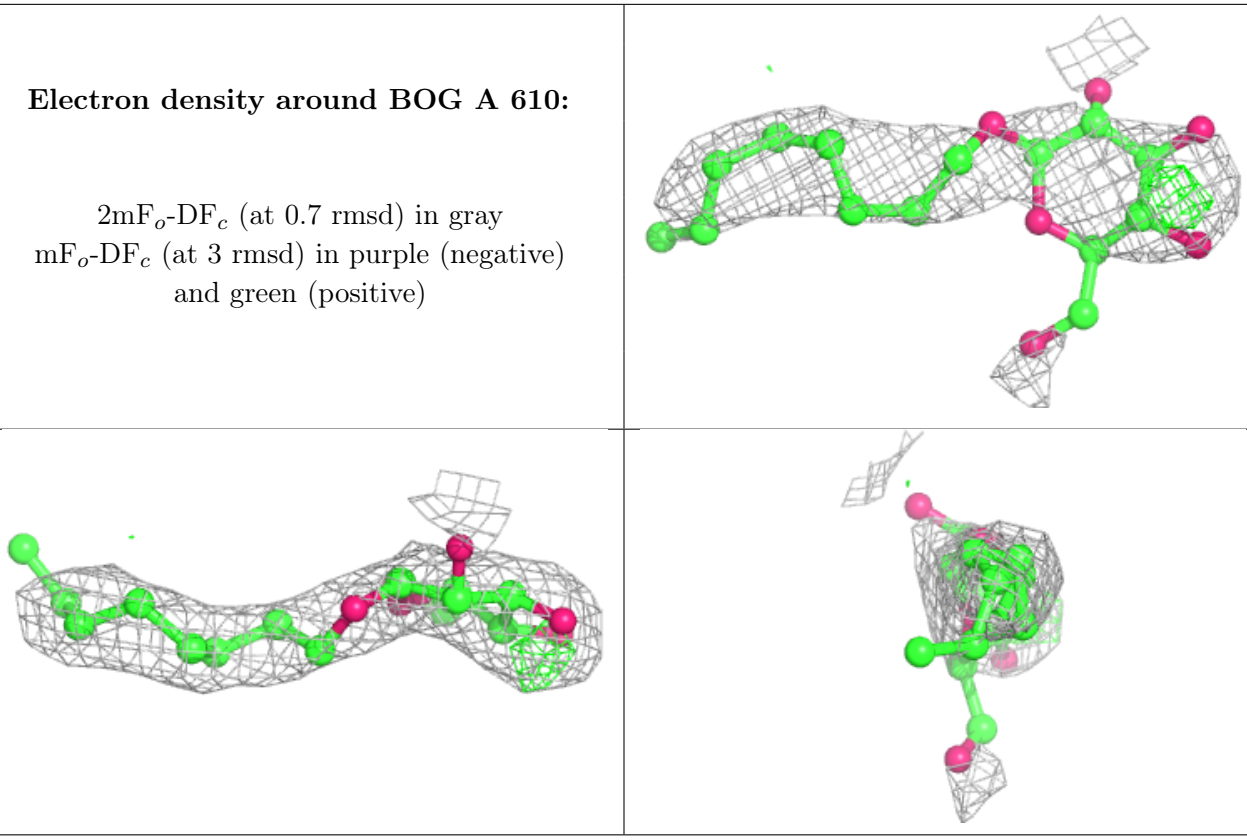
Electron density around BOG B 602:

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



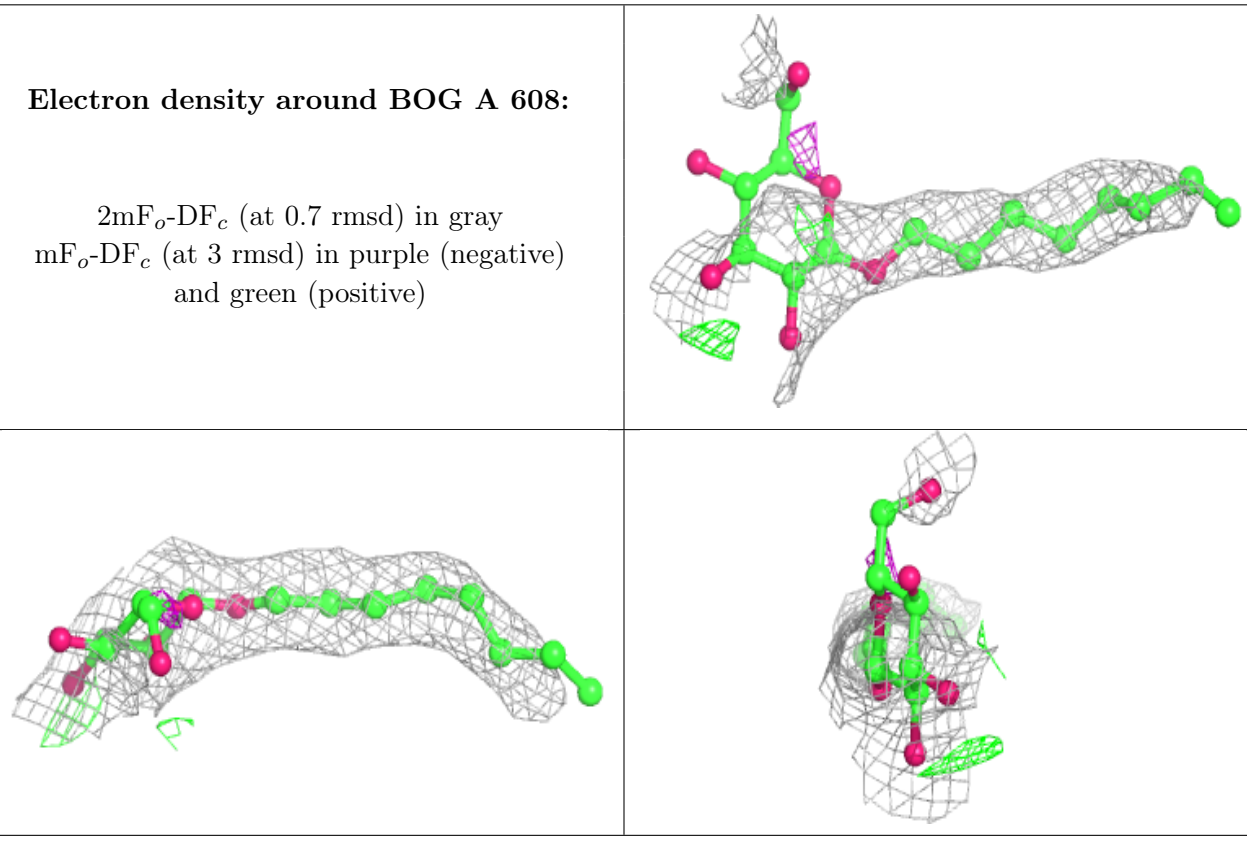
Electron density around BOG A 610:

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



Electron density around BOG A 608:

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



Electron density around BOG A 606:

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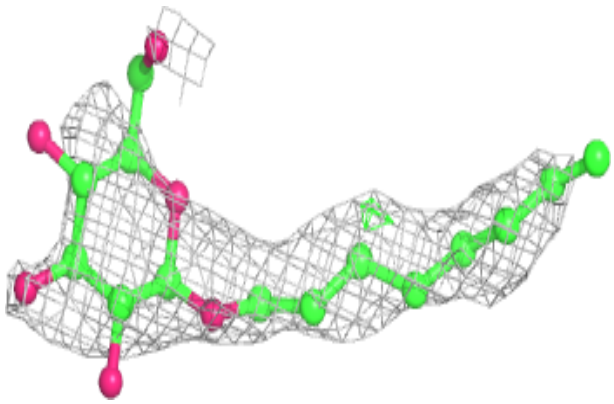
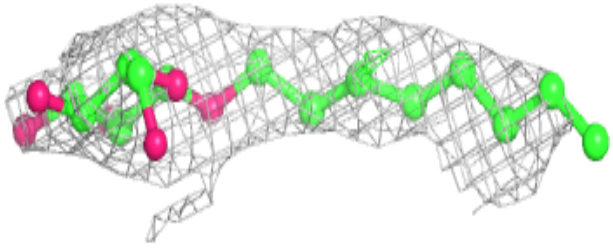
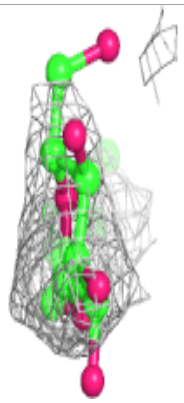
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

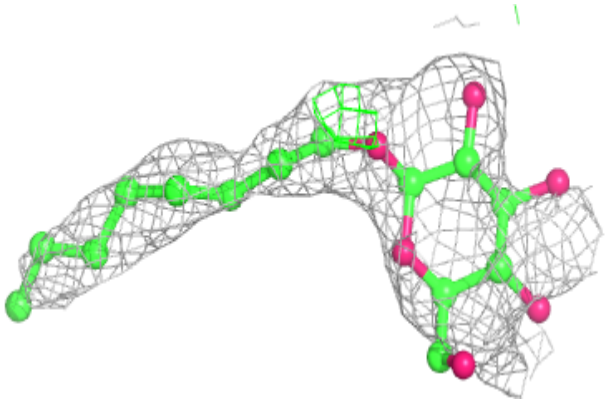


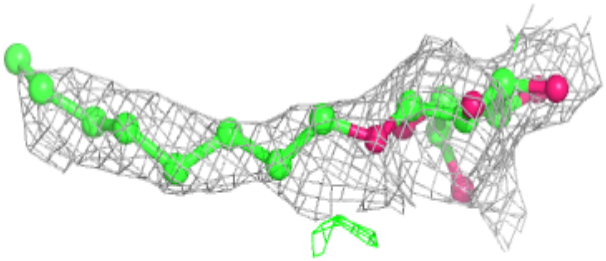


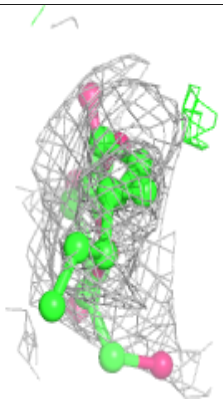


Electron density around BOG B 601:

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

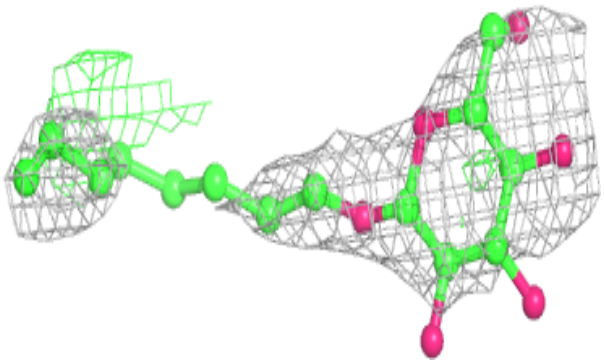


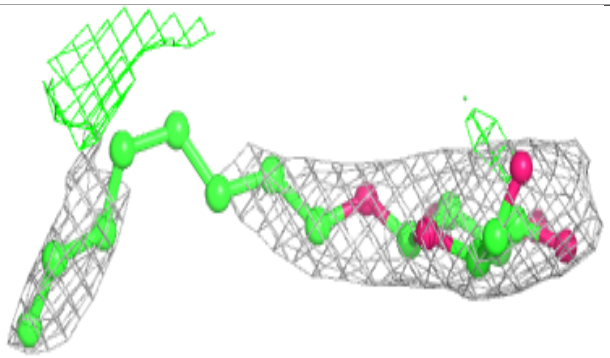


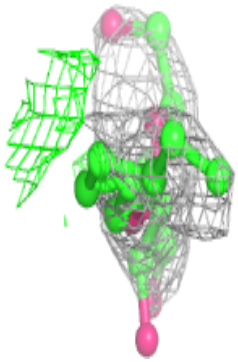


Electron density around BOG B 608:

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

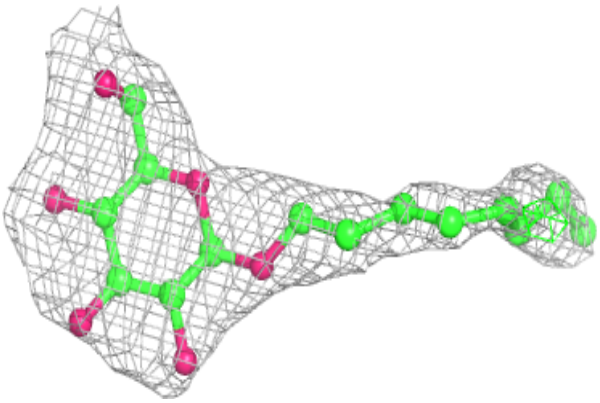


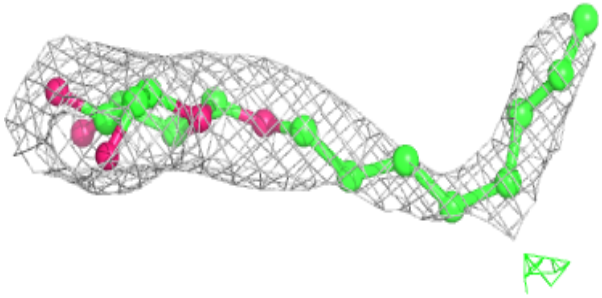


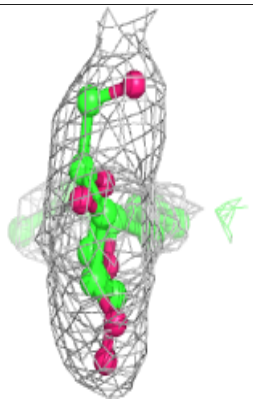


Electron density around BOG A 607:

$2mF_o$-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.