



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:59 am BST

PDB ID : 1C2T
Title : NEW INSIGHTS INTO INHIBITOR DESIGN FROM THE CRYSTAL STRUCTURE AND NMR STUDIES OF E. COLI GAR TRANSFORMYLASE IN COMPLEX WITH BETA-GAR AND 10-FORMYL-5,8,10-TRIDE AZAFOLIC ACID.
Authors : Greasley, S.E.; Yamashita, M.M.; Cai, H.; Benkovic, S.J.; Boger, D.L.; Wilson, I.A.
Deposited on : 1999-07-26
Resolution : 2.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

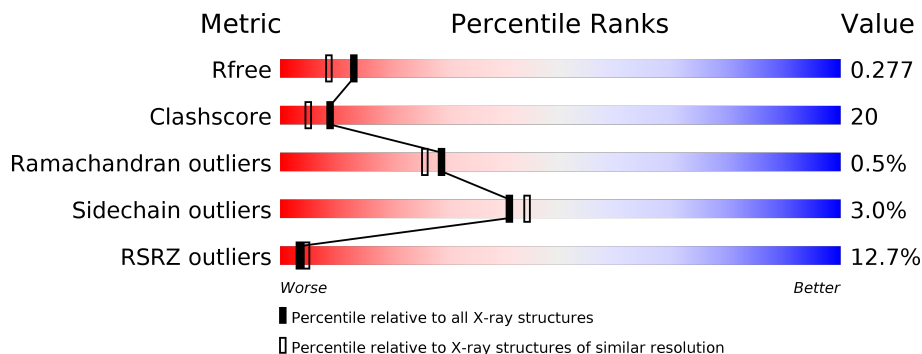
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

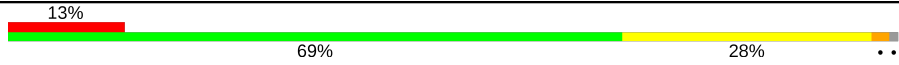

The reported resolution of this entry is 2.10 Å.

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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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

Mol	Chain	Length	Quality of chain
1	A	212	
1	B	212	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GAR	A	223	-	-	X	-
3	GAR	B	221	-	-	X	X

2 Entry composition [i](#)

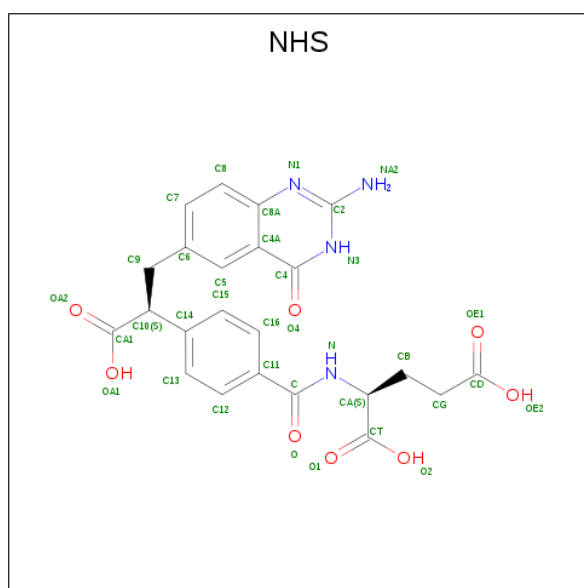
There are 4 unique types of molecules in this entry. The entry contains 3537 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLYCINAMIDE RIBONUCLEOTIDE TRANSFORMYLASE.

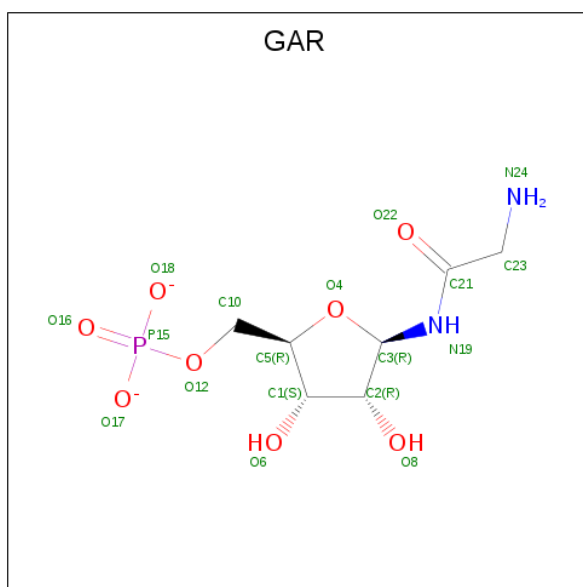
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	209	Total 1617	C 1024	N 287	O 301	S 5	0	0	0
1	B	209	Total 1617	C 1024	N 287	O 301	S 5	0	0	0

- Molecule 2 is 10-FORMYL-5,8,10-TRIDEAZAFOLIC ACID (three-letter code: NHS) (formula: $C_{23}H_{22}N_4O_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 35	C 23	N 4	O 8	0	0
2	B	1	Total 35	C 23	N 4	O 8	0	0

- Molecule 3 is GLYCINAMIDE RIBONUCLEOTIDE (three-letter code: GAR) (formula: $C_7H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	18	7	2	8	1	0	0
3	B	1	18	7	2	8	1	0	0

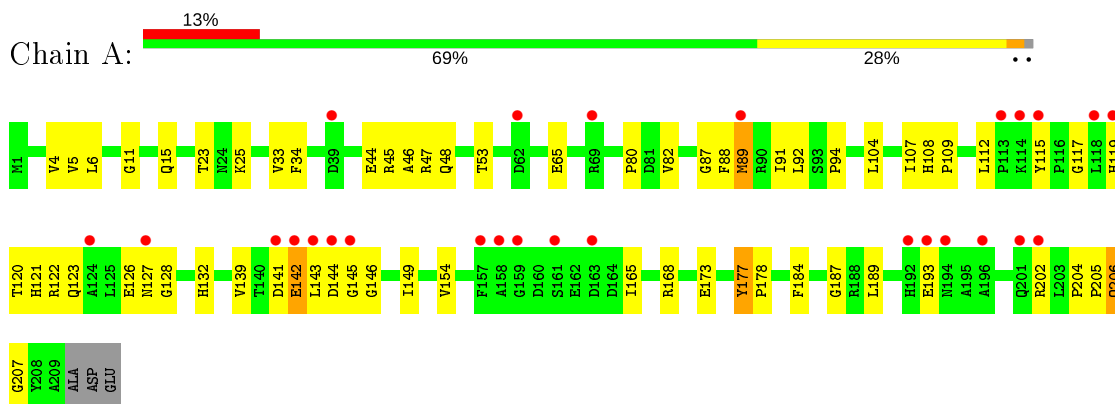
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	94	94	94	0	0
4	B	103	103	103	0	0

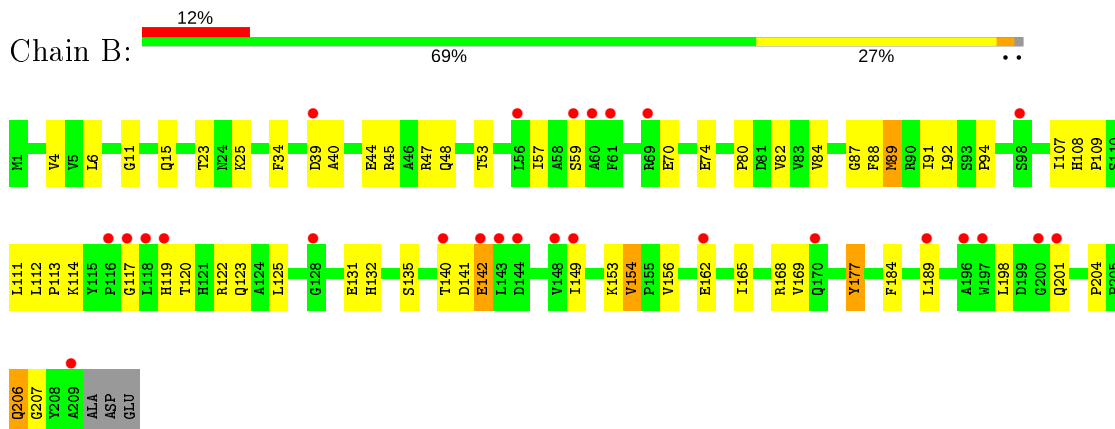
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: GLYCINAMIDE RIBONUCLEOTIDE TRANSFORMYLASE



- Molecule 1: GLYCINAMIDE RIBONUCLEOTIDE TRANSFORMYLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.25Å 112.79Å 46.96Å 90.00° 101.50° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 17.46 – 2.10	Depositor EDS
% Data completeness (in resolution range)	78.9 (20.00-2.10) 92.0 (17.46-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.11Å)	Xtrriage
Refinement program	CNS 0.3	Depositor
R, R_{free}	0.227 , 0.265 0.239 , 0.277	Depositor DCC
R_{free} test set	1085 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtrriage
Anisotropy	0.606	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3537	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.93% 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

5.1 Standard geometry

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

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	0.50	0/1654	0.73	0/2248
1	B	0.45	0/1654	0.68	0/2248
All	All	0.48	0/3308	0.70	0/4496

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1617	0	1592	74	0
1	B	1617	0	1592	56	0
2	A	35	0	20	5	0
2	B	35	0	21	5	0
3	A	18	0	13	8	0
3	B	18	0	13	7	0
4	A	94	0	0	5	0
4	B	103	0	0	3	0
All	All	3537	0	3251	134	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 20.

All (134) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:MET:HE3	3:B:221:GAR:H241	1.04	1.13
1:B:89:MET:CE	3:B:221:GAR:H241	1.64	1.09
1:A:91:ILE:HD12	1:A:143:LEU:HD11	1.42	1.02
1:B:89:MET:HE3	3:B:221:GAR:N24	1.76	1.00
1:A:202:ARG:NH2	1:B:113:PRO:HG3	1.76	0.99
1:A:202:ARG:CZ	1:B:113:PRO:HG3	1.95	0.96
1:B:140:THR:HG22	1:B:142:GLU:HG2	1.52	0.92
2:A:222:NHS:H10	3:A:223:GAR:N24	1.89	0.86
1:B:87:GLY:HA2	3:B:221:GAR:H19	1.41	0.84
1:A:112:LEU:HD21	1:A:144:ASP:HB3	1.59	0.83
1:A:89:MET:CE	3:A:223:GAR:H241	1.91	0.83
1:B:131:GLU:OE2	1:B:153:LYS:HD3	1.80	0.81
1:A:91:ILE:CD1	1:A:143:LEU:HD11	2.11	0.81
1:A:154:VAL:HG21	1:A:168:ARG:HG2	1.63	0.80
1:A:89:MET:CE	3:A:223:GAR:N24	2.47	0.78
1:A:23:THR:OG1	1:A:25:LYS:HG3	1.84	0.78
1:B:23:THR:OG1	1:B:25:LYS:HG3	1.83	0.77
1:B:132:HIS:HB3	1:B:165:ILE:HD11	1.67	0.77
1:A:11:GLY:HA2	4:A:227:HOH:O	1.86	0.76
1:A:109:PRO:HG2	1:A:120:THR:HG21	1.69	0.75
1:B:15:GLN:HE21	1:B:45:ARG:HH11	1.33	0.74
1:A:91:ILE:HD12	1:A:143:LEU:CD1	2.18	0.73
1:A:119:HIS:O	1:A:123:GLN:HG3	1.90	0.72
1:B:11:GLY:HA2	4:B:289:HOH:O	1.89	0.71
1:A:89:MET:HE1	3:A:223:GAR:H241	1.56	0.69
1:A:112:LEU:HD21	1:A:144:ASP:O	1.94	0.67
1:A:112:LEU:CD2	1:A:144:ASP:O	2.42	0.67
2:A:222:NHS:H10	3:A:223:GAR:C23	2.24	0.67
1:A:206:GLN:CA	1:A:206:GLN:HE21	2.07	0.67
1:A:115:TYR:CE2	1:A:123:GLN:HB3	2.28	0.67
1:A:15:GLN:HE21	1:A:45:ARG:HH11	1.41	0.66
1:B:206:GLN:HG3	1:B:206:GLN:O	1.94	0.66
1:A:132:HIS:O	1:A:154:VAL:HG12	1.97	0.65
1:B:156:VAL:HG22	1:B:165:ILE:HD12	1.79	0.65
2:A:222:NHS:H10	3:A:223:GAR:H231	1.80	0.64
1:B:23:THR:HG21	1:B:25:LYS:NZ	2.13	0.64
1:A:126:GLU:C	1:A:128:GLY:H	2.02	0.63
1:B:162:GLU:HG2	4:B:261:HOH:O	1.98	0.63
1:A:122:ARG:O	1:A:126:GLU:HG3	1.99	0.63
1:A:154:VAL:CG2	1:A:168:ARG:HG2	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:THR:HG21	1:A:25:LYS:NZ	2.14	0.62
1:A:94:PRO:HA	1:A:141:ASP:OD1	1.99	0.62
1:B:47:ARG:NH1	1:B:53:THR:OG1	2.33	0.61
1:B:119:HIS:O	1:B:123:GLN:HG3	2.00	0.61
1:B:132:HIS:CB	1:B:165:ILE:HD11	2.30	0.59
2:B:220:NHS:H10	3:B:221:GAR:H231	1.84	0.58
1:A:47:ARG:NH1	1:A:53:THR:OG1	2.36	0.58
1:A:132:HIS:HB3	1:A:165:ILE:HD11	1.85	0.57
1:B:140:THR:HG22	1:B:142:GLU:CG	2.32	0.57
1:B:4:VAL:HG23	1:B:80:PRO:HB3	1.85	0.57
1:B:6:LEU:HD22	1:B:34:PHE:HB2	1.86	0.56
1:A:108:HIS:HE1	1:A:117:GLY:O	1.89	0.55
1:A:132:HIS:CB	1:A:165:ILE:HD11	2.37	0.55
2:B:220:NHS:H10	3:B:221:GAR:C23	2.37	0.54
1:B:149:ILE:HD13	1:B:189:LEU:HD21	1.88	0.54
1:B:107:ILE:O	1:B:107:ILE:HG23	2.08	0.54
1:B:15:GLN:NE2	1:B:45:ARG:HH11	2.05	0.54
1:B:94:PRO:HA	1:B:141:ASP:OD1	2.08	0.53
1:B:141:ASP:HA	2:B:220:NHS:HN22	1.72	0.53
1:A:88:PHE:CE2	1:A:92:LEU:HD21	2.44	0.53
1:A:206:GLN:HB2	4:A:280:HOH:O	2.09	0.53
1:B:82:VAL:HB	1:B:184:PHE:CE2	2.43	0.53
1:A:107:ILE:O	1:A:107:ILE:HG23	2.07	0.53
1:A:88:PHE:HE2	1:A:92:LEU:HD21	1.73	0.53
1:A:206:GLN:HA	1:A:206:GLN:HE21	1.72	0.52
1:A:115:TYR:CE2	1:A:123:GLN:CB	2.91	0.52
1:A:89:MET:HA	1:A:89:MET:HE2	1.91	0.52
1:A:112:LEU:HD23	1:A:144:ASP:O	2.11	0.50
2:B:220:NHS:H10	3:B:221:GAR:N24	2.27	0.50
1:B:198:LEU:O	1:B:201:GLN:HG2	2.13	0.49
1:A:82:VAL:HB	1:A:184:PHE:CE2	2.48	0.49
1:A:15:GLN:NE2	1:A:45:ARG:HH11	2.10	0.48
1:B:88:PHE:HE2	1:B:92:LEU:HD21	1.78	0.48
1:A:187:GLY:HA2	4:A:278:HOH:O	2.13	0.48
1:A:204:PRO:HG2	1:A:207:GLY:O	2.12	0.48
1:A:126:GLU:C	1:A:128:GLY:N	2.67	0.48
1:A:126:GLU:O	1:A:128:GLY:N	2.46	0.48
1:B:109:PRO:HG2	1:B:120:THR:HG21	1.96	0.48
1:A:65:GLU:HA	4:A:249:HOH:O	2.13	0.48
1:B:88:PHE:CE2	1:B:92:LEU:HD21	2.49	0.47
1:A:142:GLU:HB2	1:A:145:GLY:HA3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ILE:HD13	1:A:189:LEU:HD21	1.96	0.47
1:A:132:HIS:O	1:A:154:VAL:N	2.48	0.47
1:A:23:THR:HG21	1:A:25:LYS:HZ1	1.78	0.47
1:B:112:LEU:HD12	1:B:135:SER:HB2	1.96	0.47
1:B:156:VAL:HG22	1:B:165:ILE:CD1	2.44	0.47
1:B:154:VAL:HG13	1:B:168:ARG:HD2	1.96	0.47
1:B:140:THR:CG2	1:B:142:GLU:HG2	2.34	0.46
1:B:132:HIS:O	1:B:154:VAL:N	2.48	0.46
1:B:57:ILE:HG22	1:B:59:SER:H	1.80	0.46
1:A:89:MET:HA	1:A:89:MET:CE	2.45	0.46
1:A:109:PRO:HG2	1:A:120:THR:CG2	2.44	0.45
1:A:120:THR:OG1	1:A:121:HIS:N	2.49	0.45
1:A:132:HIS:C	1:A:154:VAL:HG12	2.36	0.45
1:B:108:HIS:HE1	1:B:117:GLY:O	2.00	0.45
1:B:114:LYS:HG3	4:B:266:HOH:O	2.16	0.45
1:B:70:GLU:O	1:B:74:GLU:HG2	2.17	0.45
1:A:6:LEU:HD22	1:A:34:PHE:HB2	1.99	0.45
1:B:165:ILE:O	1:B:169:VAL:HG23	2.17	0.45
1:B:204:PRO:HG2	1:B:207:GLY:O	2.17	0.44
1:A:177:TYR:HB3	1:A:178:PRO:HD3	2.00	0.44
1:A:89:MET:HE3	3:A:223:GAR:N24	2.29	0.44
1:B:89:MET:CE	1:B:89:MET:HA	2.47	0.44
1:A:104:LEU:HG	1:A:139:VAL:HB	1.99	0.44
1:B:122:ARG:HH11	1:B:122:ARG:HG3	1.83	0.44
1:A:142:GLU:HG2	1:A:142:GLU:H	1.34	0.44
1:B:23:THR:HG21	1:B:25:LYS:HZ1	1.80	0.44
1:A:144:ASP:OD1	2:A:222:NHS:OA2	2.36	0.44
1:B:84:VAL:HG13	1:B:177:TYR:OH	2.18	0.43
1:A:5:VAL:C	1:A:6:LEU:HD23	2.39	0.43
1:B:39:ASP:O	1:B:40:ALA:C	2.56	0.43
1:B:91:ILE:HA	2:B:220:NHS:C8	2.49	0.43
1:B:4:VAL:CG2	1:B:80:PRO:HB3	2.48	0.43
1:B:142:GLU:HG2	1:B:142:GLU:H	1.24	0.42
1:B:122:ARG:HG3	1:B:122:ARG:NH1	2.33	0.42
1:B:125:LEU:HD22	1:B:162:GLU:CD	2.39	0.42
1:A:115:TYR:CD2	1:A:123:GLN:NE2	2.88	0.42
1:B:125:LEU:HB3	1:B:162:GLU:OE2	2.19	0.42
1:A:44:GLU:O	1:A:48:GLN:HG2	2.20	0.42
1:A:112:LEU:HD12	1:A:112:LEU:HA	1.83	0.42
1:A:87:GLY:HA2	3:A:223:GAR:H19	1.84	0.42
1:A:115:TYR:HD2	1:A:123:GLN:NE2	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:THR:CB	1:A:25:LYS:HZ2	2.31	0.41
1:A:193:GLU:HB2	4:A:313:HOH:O	2.19	0.41
1:A:141:ASP:HA	2:A:222:NHS:HN22	1.86	0.41
1:A:4:VAL:HG23	1:A:80:PRO:HB3	2.02	0.41
1:B:44:GLU:O	1:B:48:GLN:HG2	2.21	0.41
1:B:111:LEU:O	1:B:114:LYS:HB2	2.20	0.41
1:A:143:LEU:HA	1:A:143:LEU:HD12	1.57	0.41
1:A:108:HIS:CD2	1:A:112:LEU:HD13	2.56	0.40
1:A:107:ILE:HD11	1:A:173:GLU:HG2	2.03	0.40
1:A:33:VAL:HG21	1:A:46:ALA:HB2	2.02	0.40
1:A:154:VAL:HG22	1:A:168:ARG:CD	2.51	0.40
1:A:142:GLU:O	1:A:145:GLY:N	2.36	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	207/212 (98%)	194 (94%)	11 (5%)	2 (1%)	15	11
1	B	207/212 (98%)	200 (97%)	7 (3%)	0	100	100
All	All	414/424 (98%)	394 (95%)	18 (4%)	2 (0%)	29	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	ASN
1	A	146	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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	169/171 (99%)	164 (97%)	5 (3%)	41	44
1	B	169/171 (99%)	164 (97%)	5 (3%)	41	44
All	All	338/342 (99%)	328 (97%)	10 (3%)	41	44

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	MET
1	A	142	GLU
1	A	177	TYR
1	A	205	PRO
1	A	206	GLN
1	B	89	MET
1	B	142	GLU
1	B	154	VAL
1	B	177	TYR
1	B	206	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	99	HIS
1	A	123	GLN
1	A	127	ASN
1	A	206	GLN
1	B	15	GLN
1	B	123	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	NHS	B	220	-	27,37,37	2.76	14 (51%)	35,52,52	2.88	9 (25%)
2	NHS	A	222	-	27,37,37	2.53	13 (48%)	35,52,52	2.73	11 (31%)
3	GAR	B	221	-	17,18,18	1.57	3 (17%)	21,26,26	1.44	4 (19%)
3	GAR	A	223	-	17,18,18	1.32	3 (17%)	21,26,26	1.47	3 (14%)

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	NHS	B	220	-	-	7/19/29/29	0/3/3/3
2	NHS	A	222	-	-	8/19/29/29	0/3/3/3
3	GAR	B	221	-	-	8/12/28/28	0/1/1/1
3	GAR	A	223	-	-	6/12/28/28	0/1/1/1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	220	NHS	C8A-N1	6.41	1.48	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	222	NHS	C2-N3	6.33	1.46	1.35
2	B	220	NHS	C2-N3	5.96	1.46	1.35
2	A	222	NHS	C8A-N1	5.07	1.45	1.37
3	B	221	GAR	C3-N19	4.06	1.48	1.43
2	B	220	NHS	CA-N	3.87	1.51	1.46
2	B	220	NHS	C13-C14	3.77	1.45	1.39
2	A	222	NHS	C15-C14	3.48	1.44	1.39
2	A	222	NHS	C14-C10	3.41	1.62	1.53
3	B	221	GAR	P15-O18	-3.26	1.42	1.54
2	B	220	NHS	C13-C12	3.20	1.44	1.38
2	B	220	NHS	C12-C11	3.17	1.44	1.39
3	A	223	GAR	P15-O18	-3.07	1.43	1.54
2	B	220	NHS	C4-N3	3.01	1.38	1.33
2	A	222	NHS	C9-C10	2.97	1.60	1.55
2	A	222	NHS	C4-N3	2.93	1.38	1.33
2	B	220	NHS	C8-C7	2.92	1.42	1.36
2	B	220	NHS	C-N	2.91	1.40	1.34
2	A	222	NHS	C4A-C8A	2.90	1.47	1.41
2	B	220	NHS	CB-CA	2.89	1.57	1.53
2	A	222	NHS	CA-N	2.83	1.50	1.46
2	A	222	NHS	C13-C14	2.64	1.43	1.39
2	B	220	NHS	C4A-C8A	2.50	1.46	1.41
3	B	221	GAR	P15-O16	2.49	1.58	1.50
2	A	222	NHS	C8-C7	2.44	1.41	1.36
3	A	223	GAR	P15-O16	2.39	1.58	1.50
2	B	220	NHS	C14-C10	2.36	1.59	1.53
2	A	222	NHS	C12-C11	2.30	1.43	1.39
2	B	220	NHS	C15-C14	2.25	1.42	1.39
2	A	222	NHS	C16-C11	2.20	1.43	1.39
2	A	222	NHS	CB-CA	2.14	1.56	1.53
3	A	223	GAR	P15-O12	-2.11	1.53	1.60
2	B	220	NHS	C11-C	2.01	1.54	1.50

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	222	NHS	C4A-C4-N3	-9.74	117.61	124.40
2	B	220	NHS	C4A-C4-N3	-9.46	117.80	124.40
2	B	220	NHS	CA-N-C	7.48	131.98	122.34
2	B	220	NHS	C4A-C8A-N1	-6.35	120.15	123.60
2	A	222	NHS	C4A-C8A-N1	-6.03	120.32	123.60
2	A	222	NHS	C6-C9-C10	5.35	123.24	113.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	222	NHS	CA-N-C	4.74	128.46	122.34
2	B	220	NHS	C6-C9-C10	4.63	121.91	113.31
2	B	220	NHS	C4-N3-C2	4.41	122.93	115.93
2	B	220	NHS	N1-C2-N3	-4.34	121.43	127.22
2	A	222	NHS	C4-N3-C2	4.07	122.39	115.93
2	A	222	NHS	N1-C2-N3	-4.05	121.82	127.22
2	B	220	NHS	C9-C10-CA1	3.60	118.01	112.00
2	A	222	NHS	C9-C10-CA1	3.49	117.81	112.00
3	A	223	GAR	O12-P15-O16	-3.23	97.40	106.47
3	A	223	GAR	C3-N19-C21	3.12	127.96	122.45
2	A	222	NHS	CG-CB-CA	2.86	118.83	113.04
3	B	221	GAR	C3-N19-C21	2.86	127.49	122.45
2	B	220	NHS	CG-CB-CA	2.82	118.74	113.04
3	B	221	GAR	O12-P15-O16	-2.64	99.06	106.47
3	B	221	GAR	C1-C2-C3	2.53	106.23	101.42
2	A	222	NHS	C4-C4A-C8A	-2.53	116.39	118.59
3	B	221	GAR	O18-P15-O12	2.45	113.26	106.73
3	A	223	GAR	O18-P15-O12	2.24	112.69	106.73
2	B	220	NHS	C4-C4A-C8A	-2.11	116.75	118.59
2	A	222	NHS	C6-C5-C4A	-2.08	119.53	122.65
2	A	222	NHS	C2-N1-C8A	2.04	121.37	116.33

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	220	NHS	CA1-C10-C9-C6
2	B	220	NHS	CB-CA-N-C
2	B	220	NHS	CT-CA-CB-CG
2	B	220	NHS	CA-CB-CG-CD
2	A	222	NHS	CA1-C10-C9-C6
2	A	222	NHS	CB-CA-N-C
2	A	222	NHS	CT-CA-CB-CG
2	A	222	NHS	CA-CB-CG-CD
3	B	221	GAR	O4-C3-N19-C21
3	B	221	GAR	C10-O12-P15-O16
3	B	221	GAR	C10-O12-P15-O17
3	B	221	GAR	C10-O12-P15-O18
3	B	221	GAR	N19-C21-C23-N24
3	A	223	GAR	O12-C10-C5-O4
3	A	223	GAR	C10-O12-P15-O16
3	A	223	GAR	C10-O12-P15-O17

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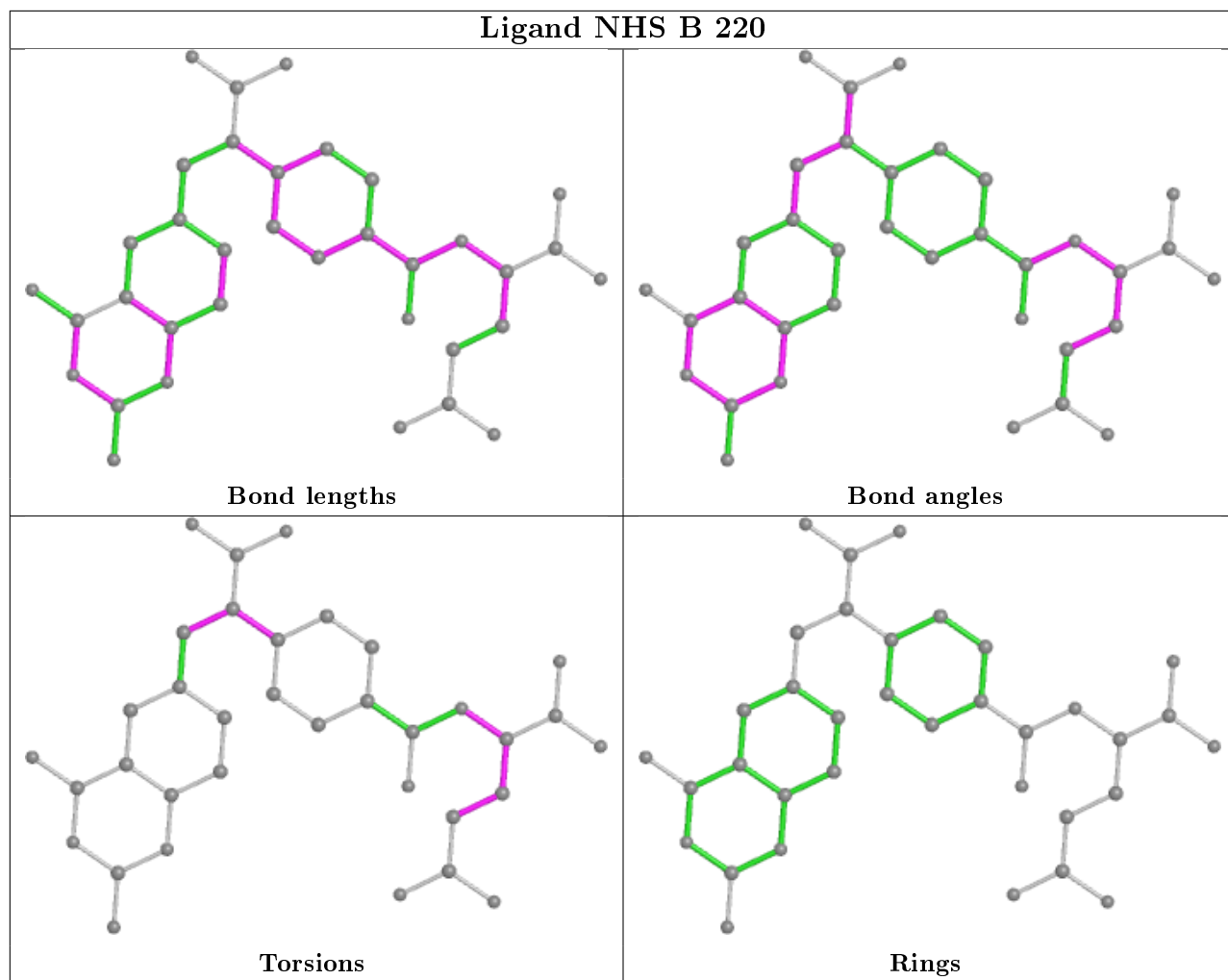
Mol	Chain	Res	Type	Atoms
3	A	223	GAR	C10-O12-P15-O18
3	B	221	GAR	O12-C10-C5-O4
3	B	221	GAR	O22-C21-C23-N24
2	A	222	NHS	CA1-C10-C14-C15
2	A	222	NHS	CA1-C10-C14-C13
3	A	223	GAR	O22-C21-C23-N24
3	B	221	GAR	O12-C10-C5-C1
3	A	223	GAR	N19-C21-C23-N24
2	B	220	NHS	N-CA-CB-CG
2	B	220	NHS	CA1-C10-C14-C13
2	B	220	NHS	CA1-C10-C14-C15
2	A	222	NHS	N-CA-CB-CG
2	A	222	NHS	C14-C10-C9-C6

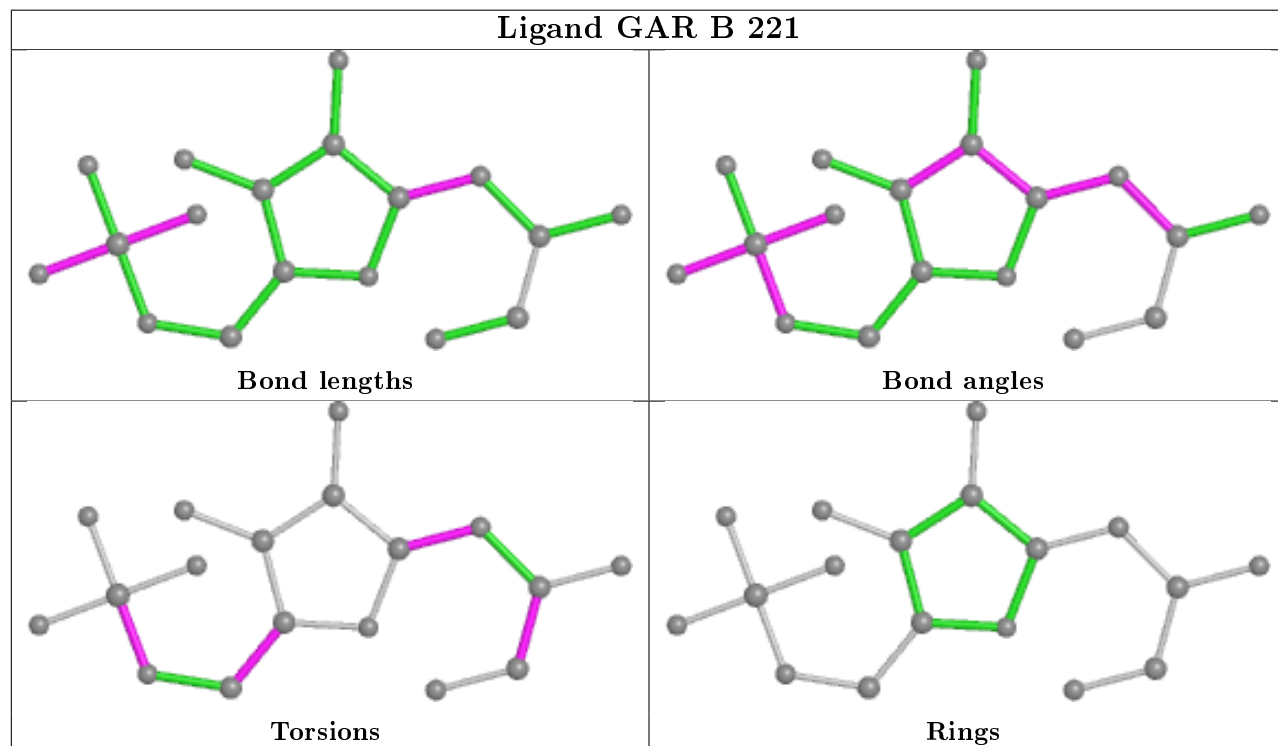
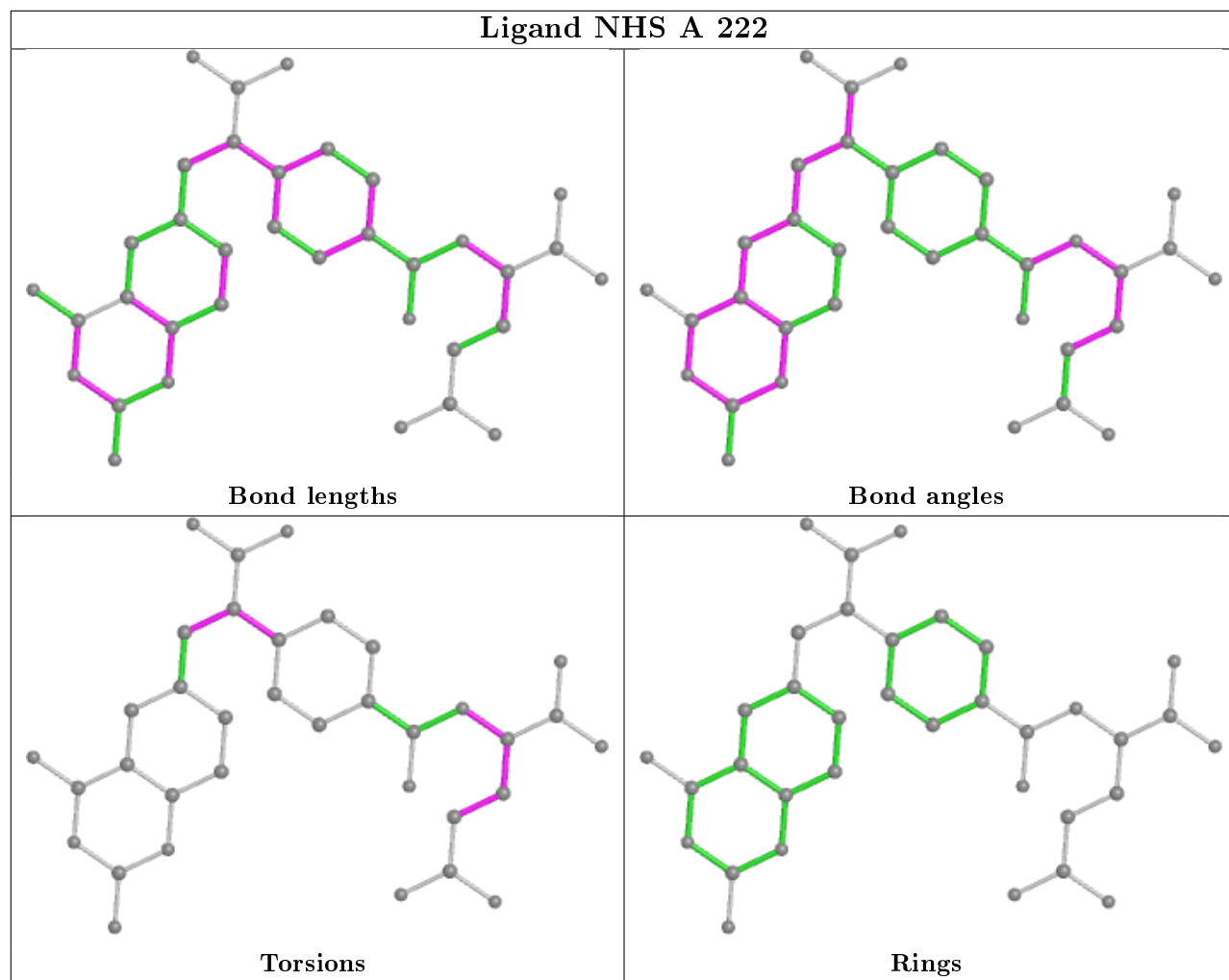
There are no ring outliers.

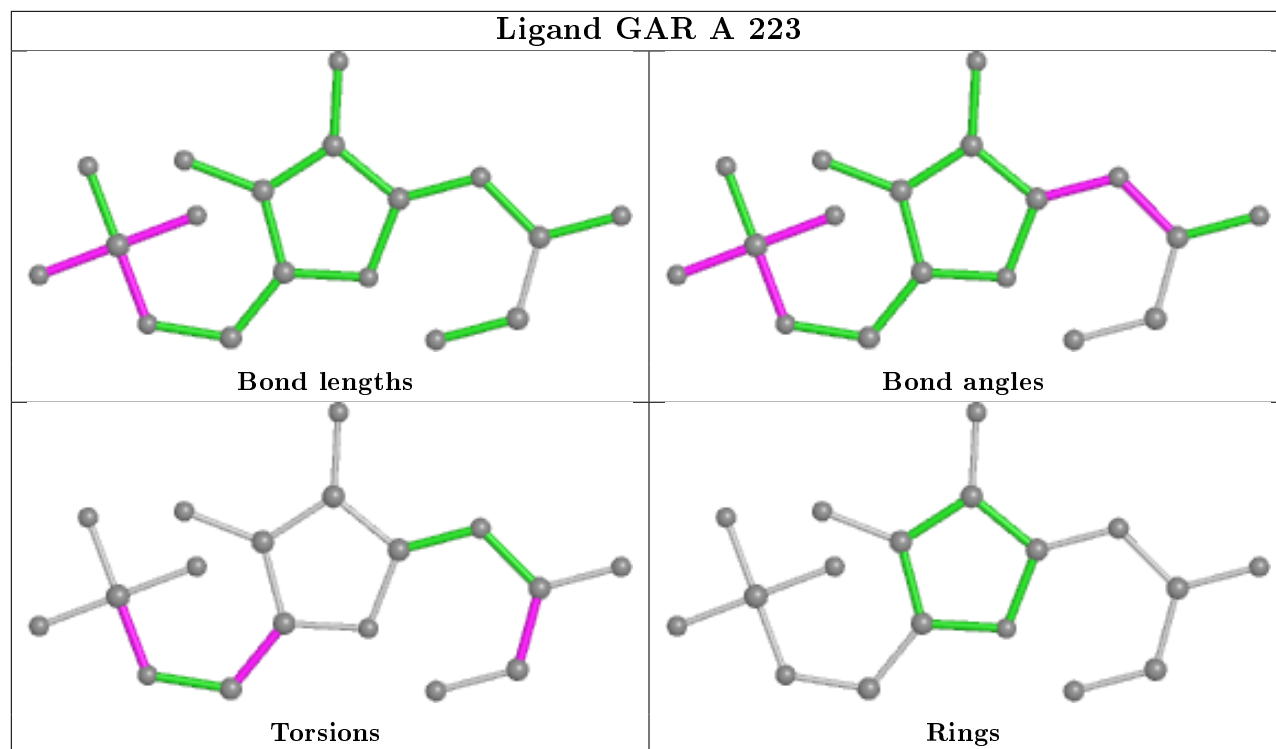
4 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	220	NHS	5	0
2	A	222	NHS	5	0
3	B	221	GAR	7	0
3	A	223	GAR	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	209/212 (98%)	0.70	27 (12%) 3 4	14, 26, 49, 52	0
1	B	209/212 (98%)	0.81	26 (12%) 4 5	16, 31, 49, 56	0
All	All	418/424 (98%)	0.76	53 (12%) 3 5	14, 30, 49, 56	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	200	GLY	7.4
1	B	39	ASP	5.4
1	B	201	GLN	4.7
1	B	209	ALA	4.3
1	B	59	SER	4.2
1	A	158	ALA	3.9
1	A	144	ASP	3.8
1	A	143	LEU	3.7
1	A	193	GLU	3.3
1	A	141	ASP	3.2
1	A	39	ASP	3.2
1	A	127	ASN	3.1
1	A	192	HIS	3.1
1	B	69	ARG	3.0
1	A	142	GLU	3.0
1	B	142	GLU	3.0
1	A	115	TYR	3.0
1	A	145	GLY	2.9
1	A	201	GLN	2.8
1	B	117	GLY	2.8
1	A	194	ASN	2.8
1	B	149	ILE	2.7
1	A	159	GLY	2.7
1	A	118	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	143	LEU	2.6
1	B	144	ASP	2.5
1	B	170	GLN	2.6
1	B	118	LEU	2.5
1	A	114	LYS	2.5
1	A	196	ALA	2.4
1	B	60	ALA	2.4
1	A	119	HIS	2.4
1	B	197	TRP	2.4
1	B	128	GLY	2.4
1	A	163	ASP	2.4
1	A	113	PRO	2.4
1	B	61	PHE	2.3
1	B	116	PRO	2.3
1	A	69	ARG	2.3
1	B	98	SER	2.2
1	A	62	ASP	2.2
1	A	89	MET	2.2
1	B	189	LEU	2.2
1	A	161	SER	2.2
1	B	196	ALA	2.2
1	A	124	ALA	2.2
1	B	162	GLU	2.1
1	B	119	HIS	2.1
1	A	157	PHE	2.1
1	A	202	ARG	2.1
1	B	56	LEU	2.0
1	B	140	THR	2.0
1	B	148	VAL	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](#)

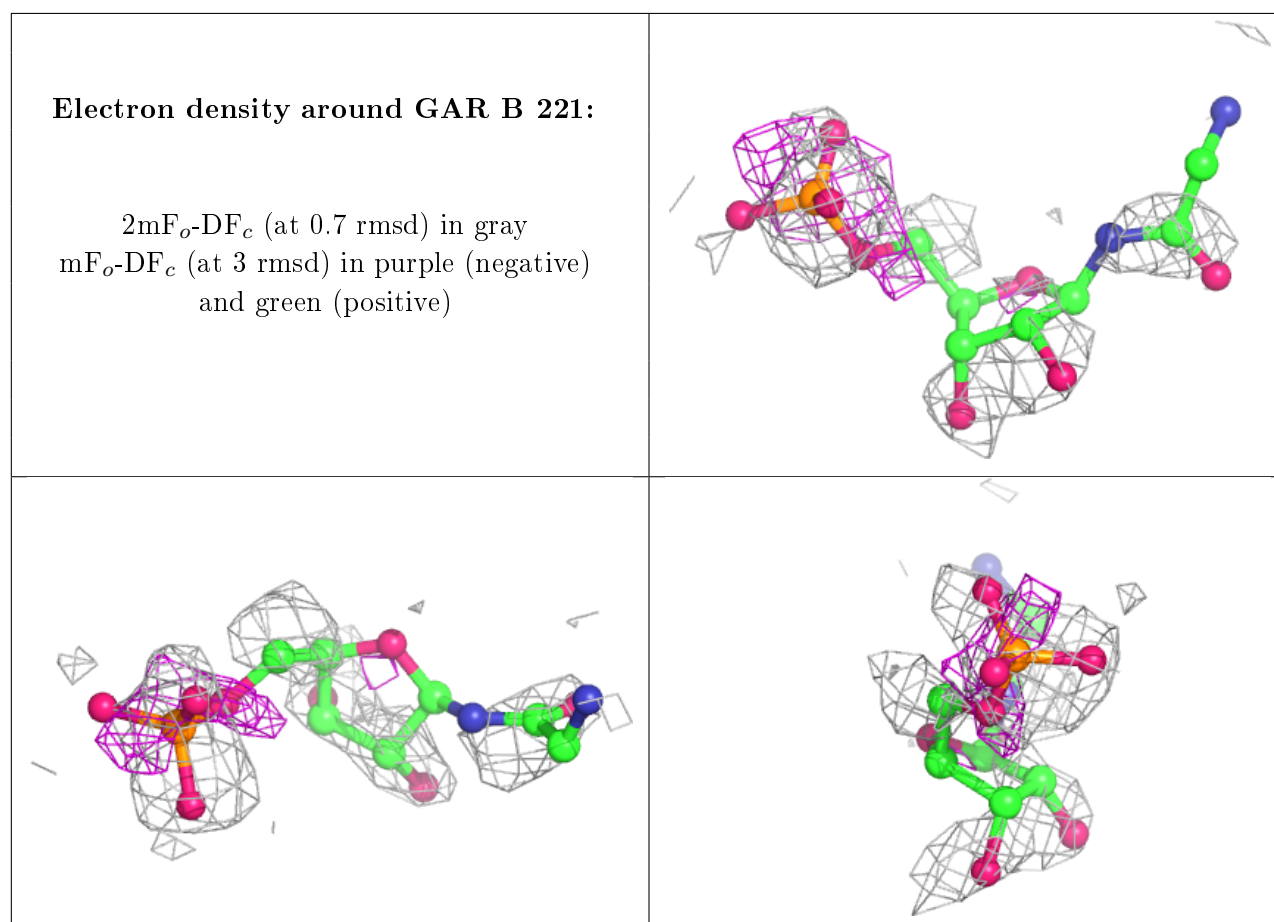
There are no carbohydrates in this entry.

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.

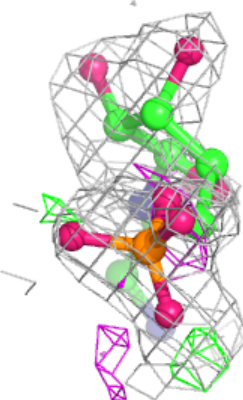
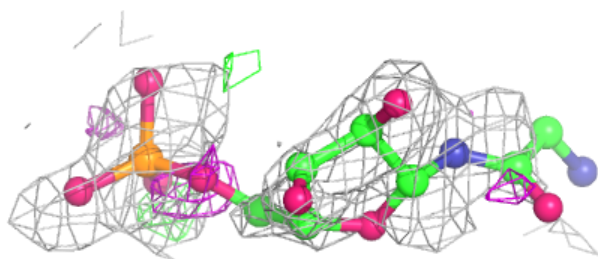
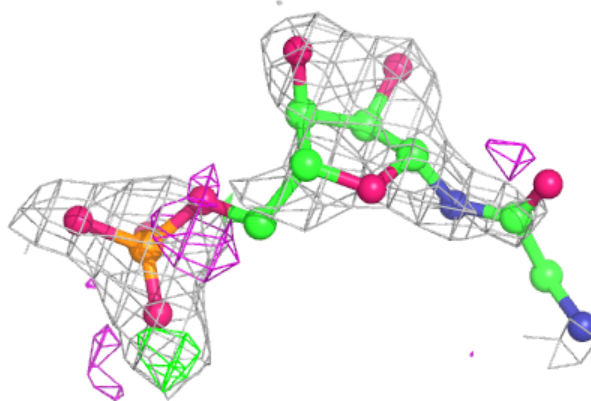
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GAR	B	221	18/18	0.64	0.49	58,61,62,62	0
3	GAR	A	223	18/18	0.70	0.39	50,56,61,61	0
2	NHS	B	220	35/35	0.77	0.23	41,47,52,53	0
2	NHS	A	222	35/35	0.79	0.20	32,43,53,53	0

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.

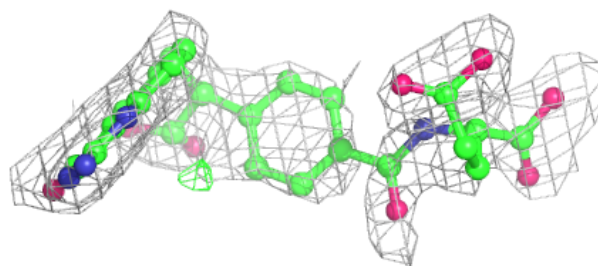
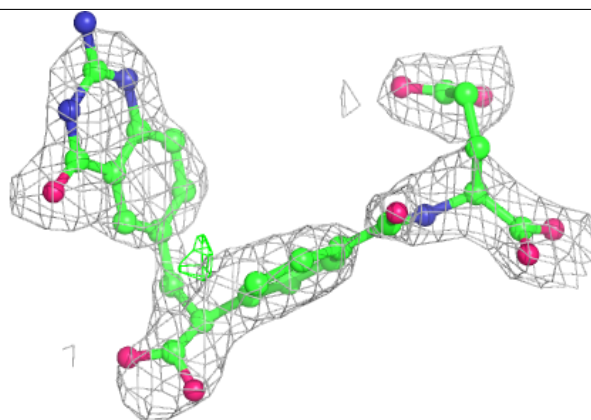


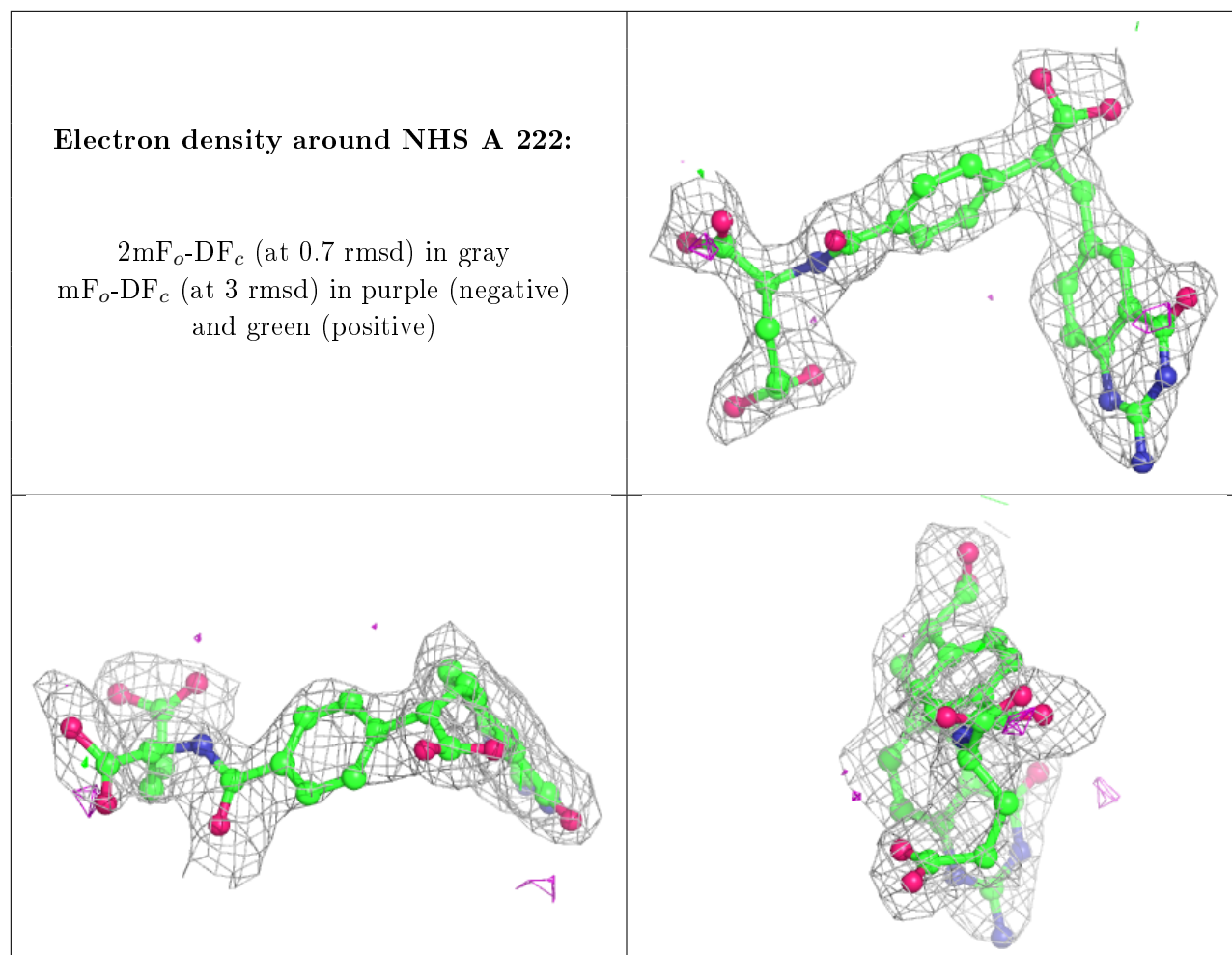
Electron density around GAR A 223:

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

**Electron density around NHS B 220:**

$2mF_o-DF_c$ (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.