



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 04:11 am BST

PDB ID : 3K34
Title : Human carbonic anhydrase II with a sulfonamide inhibitor
Authors : Behnke, C.A.; Le Trong, I.; Merritt, E.A.; Teller, D.C.; Stenkamp, R.E.
Deposited on : 2009-10-01
Resolution : 0.90 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

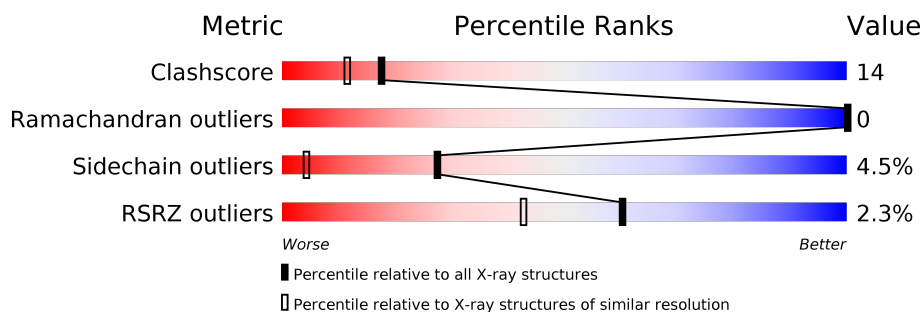
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 0.90 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))
Clashscore	141614	1132 (1.04-0.76)
Ramachandran outliers	138981	1055 (1.04-0.76)
Sidechain outliers	138945	1056 (1.04-0.76)
RSRZ outliers	127900	1028 (1.04-0.76)

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	260	

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	HGB	A	265[B]	-	X	-	-
5	GOL	A	1004	-	X	X	-
5	GOL	A	1005	-	-	X	-
5	GOL	A	1006	-	X	-	-

2 Entry composition [i](#)

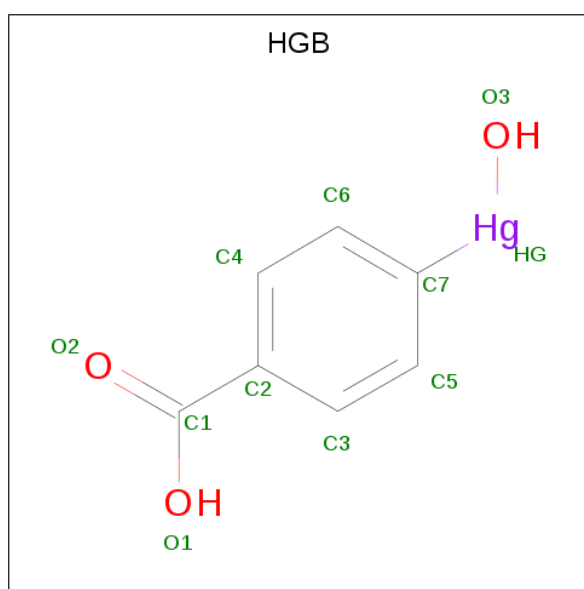
There are 6 unique types of molecules in this entry. The entry contains 2387 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Carbonic anhydrase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	10	0
			2086	1337	359	387	3			

- Molecule 2 is 4-(HYDROXYMERCURY)BENZOIC ACID (three-letter code: HGB) (formula: $C_7H_6HgO_3$).

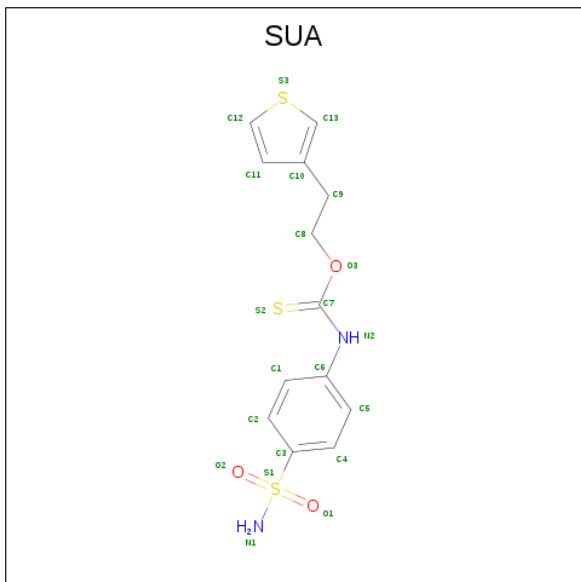


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Hg	O	0	1
			12	8	2	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

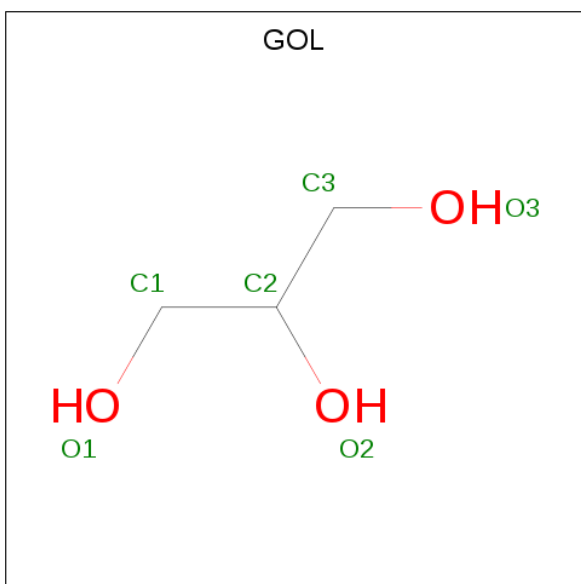
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is (4-SULFAMOYL-PHENYL)-THIOCARBAMIC ACID O-(2-THIOPHEN-3-Y L-ETHYL) ESTER (three-letter code: SUA) (formula: $C_{13}H_{14}N_2O_3S_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			21	13	2	3	3		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

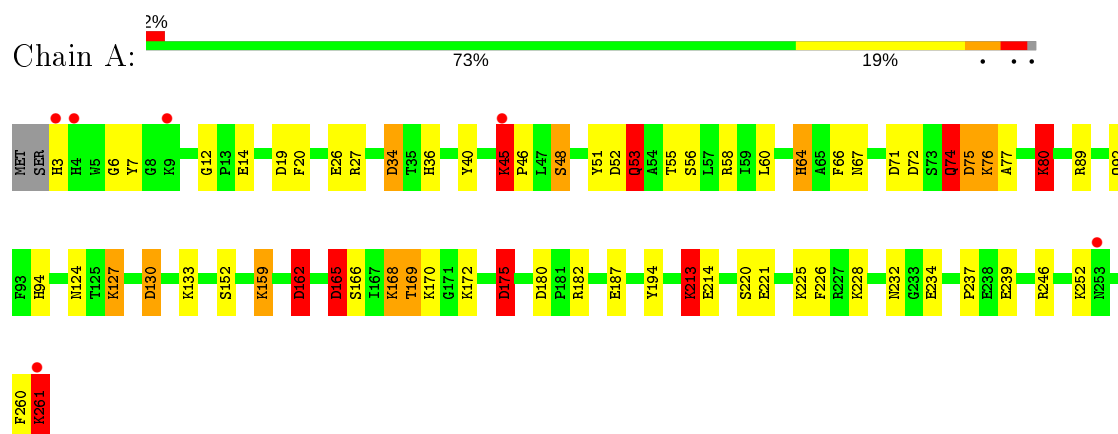
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	249	Total	O	0	0
			249	249		

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: Carbonic anhydrase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.05Å 41.27Å 71.82Å 90.00° 104.23° 90.00°	Depositor
Resolution (Å)	18.00 – 0.90 17.88 – 0.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (18.00-0.90) 88.2 (17.88-0.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 0.90Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.141 , 0.160 0.130 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	8.7	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2387	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.27% 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: GOL, ZN, HGB, SUA

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	1.51	23/2199 (1.0%)	2.07	72/2978 (2.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	14	GLU	CD-OE2	14.56	1.41	1.25
1	A	27	ARG	CD-NE	-13.62	1.23	1.46
1	A	64	HIS	CD2-NE2	-8.96	1.18	1.38
1	A	64	HIS	CE1-NE2	7.79	1.50	1.32
1	A	159	LYS	CE-NZ	7.79	1.68	1.49
1	A	27	ARG	CZ-NH2	-7.08	1.23	1.33
1	A	220[A]	SER	CB-OG	-6.64	1.33	1.42
1	A	220[B]	SER	CB-OG	-6.64	1.33	1.42
1	A	40	TYR	CE2-CZ	-6.61	1.29	1.38
1	A	234	GLU	CG-CD	-6.53	1.42	1.51
1	A	130	ASP	CA-CB	6.28	1.67	1.53
1	A	234	GLU	CD-OE1	6.25	1.32	1.25
1	A	3	HIS	C-O	6.16	1.35	1.23
1	A	226	PHE	CG-CD2	5.85	1.47	1.38
1	A	56	SER	CB-OG	5.83	1.49	1.42
1	A	152[A]	SER	CA-CB	5.68	1.61	1.52
1	A	152[B]	SER	CA-CB	5.68	1.61	1.52
1	A	232	ASN	C-N	5.53	1.43	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	221	GLU	CD-OE1	-5.23	1.19	1.25
1	A	214	GLU	CD-OE1	-5.20	1.20	1.25
1	A	182	ARG	NE-CZ	5.14	1.39	1.33
1	A	48	SER	CB-OG	-5.11	1.35	1.42
1	A	89	ARG	NE-CZ	-5.04	1.26	1.33

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	ASP	CB-CG-OD2	32.01	147.11	118.30
1	A	27	ARG	NE-CZ-NH2	-23.30	108.65	120.30
1	A	27	ARG	NE-CZ-NH1	20.24	130.42	120.30
1	A	182	ARG	NE-CZ-NH2	-20.13	110.23	120.30
1	A	175	ASP	CB-CG-OD1	-17.62	102.44	118.30
1	A	75	ASP	CB-CG-OD1	16.91	133.52	118.30
1	A	162	ASP	CA-CB-CG	14.85	146.06	113.40
1	A	130	ASP	CB-CG-OD1	14.60	131.44	118.30
1	A	162	ASP	CB-CG-OD1	-14.59	105.17	118.30
1	A	80	LYS	CD-CE-NZ	14.36	144.72	111.70
1	A	74	GLN	OE1-CD-NE2	13.78	153.59	121.90
1	A	58	ARG	NE-CZ-NH1	13.08	126.84	120.30
1	A	27	ARG	CG-CD-NE	12.74	138.55	111.80
1	A	226	PHE	CB-CG-CD1	12.65	129.66	120.80
1	A	182	ARG	NE-CZ-NH1	12.13	126.36	120.30
1	A	239	GLU	OE1-CD-OE2	11.96	137.66	123.30
1	A	74	GLN	CG-CD-OE1	-11.21	99.18	121.60
1	A	130	ASP	OD1-CG-OD2	-10.93	102.54	123.30
1	A	27	ARG	CD-NE-CZ	10.12	137.77	123.60
1	A	170	LYS	CD-CE-NZ	10.05	134.82	111.70
1	A	77	ALA	N-CA-CB	-9.99	96.11	110.10
1	A	58	ARG	NE-CZ-NH2	-9.94	115.33	120.30
1	A	165	ASP	CB-CG-OD2	-9.91	109.38	118.30
1	A	175	ASP	OD1-CG-OD2	9.57	141.49	123.30
1	A	19	ASP	CB-CG-OD1	8.71	126.14	118.30
1	A	3	HIS	O-C-N	8.63	136.52	122.70
1	A	175	ASP	CA-CB-CG	-8.45	94.81	113.40
1	A	162	ASP	OD1-CG-OD2	-8.39	107.36	123.30
1	A	221	GLU	OE1-CD-OE2	-8.37	113.26	123.30
1	A	226	PHE	CB-CG-CD2	-8.26	115.02	120.80
1	A	34	ASP	CB-CG-OD1	7.70	125.23	118.30
1	A	3	HIS	CA-C-O	-7.45	104.46	120.10
1	A	76	LYS	O-C-N	-7.44	110.80	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	168	LYS	CD-CE-NZ	7.36	128.62	111.70
1	A	246	ARG	NE-CZ-NH2	-7.11	116.74	120.30
1	A	75	ASP	OD1-CG-OD2	-6.92	110.15	123.30
1	A	71	ASP	CB-CG-OD2	-6.87	112.11	118.30
1	A	89	ARG	NE-CZ-NH1	6.85	123.72	120.30
1	A	26	GLU	OE1-CD-OE2	-6.71	115.25	123.30
1	A	252	LYS	CB-CG-CD	6.68	128.96	111.60
1	A	180	ASP	CB-CG-OD1	6.63	124.27	118.30
1	A	165	ASP	OD1-CG-OD2	6.60	135.84	123.30
1	A	20	PHE	CB-CG-CD2	-6.53	116.23	120.80
1	A	159	LYS	CG-CD-CE	6.34	130.92	111.90
1	A	34	ASP	OD1-CG-OD2	-6.33	111.28	123.30
1	A	64	HIS	ND1-CE1-NE2	-6.21	96.24	109.90
1	A	261	LYS	CA-CB-CG	-6.18	99.80	113.40
1	A	19	ASP	OD1-CG-OD2	-6.18	111.56	123.30
1	A	14	GLU	CG-CD-OE1	-6.08	106.14	118.30
1	A	40	TYR	CB-CG-CD1	6.03	124.62	121.00
1	A	225	LYS	CB-CG-CD	6.00	127.19	111.60
1	A	246	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	187[A]	GLU	OE1-CD-OE2	-5.74	116.41	123.30
1	A	187[B]	GLU	OE1-CD-OE2	-5.74	116.41	123.30
1	A	3	HIS	CA-CB-CG	-5.68	103.94	113.60
1	A	55	THR	CA-CB-CG2	-5.65	104.49	112.40
1	A	127[A]	LYS	CD-CE-NZ	-5.64	98.72	111.70
1	A	127[B]	LYS	CD-CE-NZ	-5.64	98.72	111.70
1	A	175	ASP	CB-CA-C	-5.51	99.38	110.40
1	A	14	GLU	CG-CD-OE2	5.47	129.25	118.30
1	A	237	PRO	N-CA-CB	5.44	109.83	103.30
1	A	213	LYS	CD-CE-NZ	-5.37	99.35	111.70
1	A	213	LYS	CG-CD-CE	-5.29	96.02	111.90
1	A	45	LYS	CD-CE-NZ	-5.26	99.60	111.70
1	A	52	ASP	CA-CB-CG	5.21	124.87	113.40
1	A	194	TYR	CB-CG-CD2	-5.19	117.89	121.00
1	A	51	TYR	CB-CG-CD2	-5.10	117.94	121.00
1	A	169	THR	CA-C-N	-5.07	106.04	117.20
1	A	53	GLN	CG-CD-OE1	5.06	131.73	121.60
1	A	239	GLU	CG-CD-OE2	-5.06	108.18	118.30
1	A	3	HIS	C-N-CA	5.03	134.28	121.70
1	A	72	ASP	CB-CG-OD1	5.02	122.82	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	34	ASP	Sidechain
1	A	64	HIS	Sidechain

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	2086	0	2031	49	4
2	A	12	0	4	0	0
3	A	1	0	0	0	0
4	A	21	0	12	5	0
5	A	18	0	18	8	2
6	A	249	0	0	12	6
All	All	2387	0	2065	58	7

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (58) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LYS:CE	1:A:159:LYS:NZ	1.68	1.54
5:A:1006:GOL:C1	5:A:1006:GOL:C2	1.95	1.45
5:A:1005:GOL:C1	5:A:1005:GOL:O1	1.69	1.39
5:A:1004:GOL:C1	5:A:1004:GOL:O1	1.77	1.30
1:A:45:LYS:NZ	1:A:46:PRO:HD2	1.53	1.23
1:A:45:LYS:NZ	1:A:45:LYS:HB3	1.56	1.14
1:A:45:LYS:CB	1:A:45:LYS:HZ2	1.67	1.06
1:A:124:ASN:HD21	1:A:127[B]:LYS:HE3	1.18	1.05
1:A:45:LYS:NZ	1:A:46:PRO:CD	2.21	1.03
1:A:45:LYS:HZ1	1:A:46:PRO:HD2	0.89	1.03
1:A:45:LYS:HZ1	1:A:46:PRO:CD	1.73	1.02
1:A:124:ASN:ND2	1:A:127[B]:LYS:HE3	1.76	1.00
1:A:45:LYS:NZ	1:A:45:LYS:CB	2.20	0.99
1:A:45:LYS:HB3	1:A:45:LYS:HZ2	0.81	0.97
1:A:130:ASP:OD2	1:A:133[B]:LYS:NZ	2.01	0.93
1:A:36:HIS:HD2	6:A:2197:HOH:O	1.57	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:LYS:HG3	1:A:261:LYS:O	1.66	0.85
1:A:124:ASN:HD21	1:A:127[B]:LYS:CE	1.94	0.80
5:A:1005:GOL:C2	5:A:1005:GOL:O1	2.32	0.78
1:A:36:HIS:CD2	6:A:2197:HOH:O	2.34	0.78
1:A:213:LYS:HE3	1:A:260:PHE:CE1	2.19	0.77
1:A:130:ASP:OD1	1:A:133[B]:LYS:NZ	2.22	0.72
1:A:162:ASP:OD1	6:A:2243:HOH:O	2.09	0.70
5:A:1006:GOL:O2	5:A:1006:GOL:C1	2.41	0.68
4:A:1003:SUA:S2	6:A:2235:HOH:O	2.53	0.66
1:A:45:LYS:CB	1:A:45:LYS:HZ3	2.09	0.65
1:A:7:TYR:O	6:A:2383:HOH:O	2.14	0.64
1:A:165:ASP:N	1:A:165:ASP:OD1	2.21	0.60
1:A:76:LYS:HE2	6:A:2247:HOH:O	2.02	0.59
1:A:162:ASP:CB	6:A:2243:HOH:O	2.51	0.58
1:A:168:LYS:NZ	1:A:228:LYS:O	2.37	0.58
1:A:213:LYS:HE3	1:A:260:PHE:CZ	2.39	0.57
1:A:48:SER:HB3	1:A:80:LYS:HG2	1.86	0.56
1:A:169:THR:OG1	1:A:172[B]:LYS:HD2	2.07	0.55
1:A:124:ASN:HD21	1:A:127[B]:LYS:NZ	2.06	0.54
5:A:1005:GOL:C1	5:A:1005:GOL:HO1	2.12	0.53
1:A:159:LYS:CD	1:A:159:LYS:NZ	2.64	0.53
1:A:162:ASP:HA	6:A:2243:HOH:O	2.08	0.52
1:A:45:LYS:HZ2	1:A:46:PRO:HD2	1.68	0.51
1:A:166:SER:O	1:A:172[A]:LYS:HE2	2.12	0.50
1:A:67:ASN:HD22	1:A:94:HIS:HB3	1.75	0.50
1:A:175:ASP:HB3	6:A:2420:HOH:O	2.12	0.48
1:A:162:ASP:CA	6:A:2243:HOH:O	2.60	0.48
1:A:130:ASP:CG	1:A:133[B]:LYS:NZ	2.67	0.48
5:A:1004:GOL:C3	5:A:1004:GOL:O1	2.63	0.46
1:A:159:LYS:HE3	6:A:2238:HOH:O	2.16	0.45
1:A:45:LYS:NZ	1:A:46:PRO:HD3	2.26	0.44
1:A:60:LEU:O	1:A:66:PHE:HA	2.18	0.44
1:A:92:GLN:HE21	4:A:1003:SUA:H92	1.83	0.44
1:A:92:GLN:HE21	4:A:1003:SUA:C9	2.30	0.43
1:A:92:GLN:HG2	4:A:1003:SUA:C9	2.49	0.42
1:A:92:GLN:HG2	4:A:1003:SUA:H91	2.01	0.42
1:A:6:GLY:O	1:A:12:GLY:HA2	2.20	0.41
5:A:1005:GOL:C3	5:A:1005:GOL:O1	2.68	0.41
1:A:168:LYS:NZ	6:A:2072:HOH:O	2.23	0.41
1:A:74:GLN:HB2	1:A:74:GLN:HE21	1.10	0.41
1:A:67:ASN:ND2	1:A:94:HIS:HB3	2.37	0.40

All (7) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1004:GOL:O1	6:A:2112:HOH:O[2_546]	1.13	1.07
5:A:1004:GOL:C1	6:A:2117:HOH:O[2_546]	1.39	0.81
1:A:130:ASP:CG	6:A:2208:HOH:O[1_565]	1.51	0.69
6:A:2084:HOH:O	6:A:2233:HOH:O[1_655]	1.65	0.55
1:A:130:ASP:OD2	6:A:2208:HOH:O[1_565]	1.66	0.54
1:A:130:ASP:CB	6:A:2208:HOH:O[1_565]	1.74	0.46
1:A:53:GLN:NE2	1:A:165:ASP:OD2[2_455]	2.09	0.11

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	266 / 260 (102%)	256 (96%)	10 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233 / 225 (104%)	223 (96%)	10 (4%)	29	5

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

Mol	Chain	Res	Type
1	A	45	LYS
1	A	53	GLN
1	A	74	GLN
1	A	75	ASP
1	A	80	LYS
1	A	162	ASP
1	A	165	ASP
1	A	175	ASP
1	A	213	LYS
1	A	261	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	53	GLN
1	A	67	ASN
1	A	74	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](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z > 2$
2	HGB	A	265[A]	1,6	5,10,11	5.03	1 (20%)	8,13,14	1.82	2 (25%)
5	GOL	A	1006	-	5,5,5	4.90	2 (40%)	5,5,5	3.10	3 (60%)
2	HGB	A	265[B]	1,6	5,1,11	5.03	1 (20%)	8,0,14	10.01	4 (50%)
5	GOL	A	1005	-	5,5,5	3.50	2 (40%)	5,5,5	2.54	3 (60%)
4	SUA	A	1003	3	21,22,22	7.32	6 (28%)	27,30,30	3.52	10 (37%)
5	GOL	A	1004	-	5,5,5	5.31	4 (80%)	5,5,5	2.71	1 (20%)

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	HGB	A	265[A]	1,6	-	0/0/4/6	0/1/1/1
5	GOL	A	1006	-	-	2/4/4/4	-
2	HGB	A	265[B]	1,6	-	0/0/0/6	0/1/0/1
5	GOL	A	1005	-	-	0/4/4/4	-
4	SUA	A	1003	3	-	4/16/16/16	0/2/2/2
5	GOL	A	1004	-	-	2/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1003	SUA	C7-N2	-27.92	1.11	1.34
4	A	1003	SUA	C13-C10	15.46	1.46	1.37
2	A	265[A]	HGB	C2-C1	11.05	1.58	1.47
2	A	265[B]	HGB	C2-C1	11.05	1.58	1.47
5	A	1006	GOL	C1-C2	10.58	1.95	1.51
5	A	1004	GOL	O1-C1	8.39	1.77	1.42
4	A	1003	SUA	O3-C7	6.95	1.41	1.33
5	A	1005	GOL	O1-C1	6.38	1.69	1.42
5	A	1004	GOL	C3-C2	-6.28	1.25	1.51
4	A	1003	SUA	C7-S2	4.46	1.74	1.66
5	A	1004	GOL	O2-C2	4.26	1.56	1.43
5	A	1005	GOL	O3-C3	3.79	1.58	1.42
5	A	1004	GOL	O3-C3	-3.50	1.27	1.42
4	A	1003	SUA	C6-N2	-3.09	1.35	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1003	SUA	C11-C12	2.16	1.46	1.37
5	A	1006	GOL	O1-C1	-2.13	1.33	1.42

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	265[B]	HGB	C3-C5-C7	22.45	165.32	121.56
2	A	265[B]	HGB	C6-C7-C5	-14.52	47.03	117.48
4	A	1003	SUA	O3-C7-S2	-9.69	118.48	125.10
4	A	1003	SUA	C8-O3-C7	-9.12	111.70	119.11
4	A	1003	SUA	O3-C7-N2	8.67	130.07	111.94
2	A	265[B]	HGB	C4-C6-C7	8.58	138.28	121.56
5	A	1004	GOL	O2-C2-C3	5.74	134.39	109.12
4	A	1003	SUA	S2-C7-N2	-5.24	111.45	123.02
5	A	1006	GOL	O2-C2-C3	4.96	130.97	109.12
5	A	1006	GOL	O2-C2-C1	-4.24	90.43	109.12
5	A	1005	GOL	O1-C1-C2	-4.08	90.62	110.20
4	A	1003	SUA	C6-N2-C7	2.80	135.00	130.00
5	A	1005	GOL	O3-C3-C2	-2.68	97.37	110.20
2	A	265[A]	HGB	C6-C4-C2	-2.61	117.38	121.13
2	A	265[B]	HGB	C6-C4-C2	-2.61	117.38	121.13
2	A	265[A]	HGB	C4-C6-C7	2.54	126.52	121.56
4	A	1003	SUA	C9-C10-C13	-2.51	122.02	127.24
4	A	1003	SUA	O3-C8-C9	-2.46	101.52	108.44
4	A	1003	SUA	C10-C13-S3	-2.44	110.72	112.29
5	A	1005	GOL	C3-C2-C1	2.38	120.97	111.70
4	A	1003	SUA	O1-S1-N1	2.24	110.69	107.36
4	A	1003	SUA	O2-S1-N1	2.12	110.51	107.36
5	A	1006	GOL	C3-C2-C1	2.06	119.73	111.70

There are no chirality outliers.

All (8) torsion outliers are listed below:

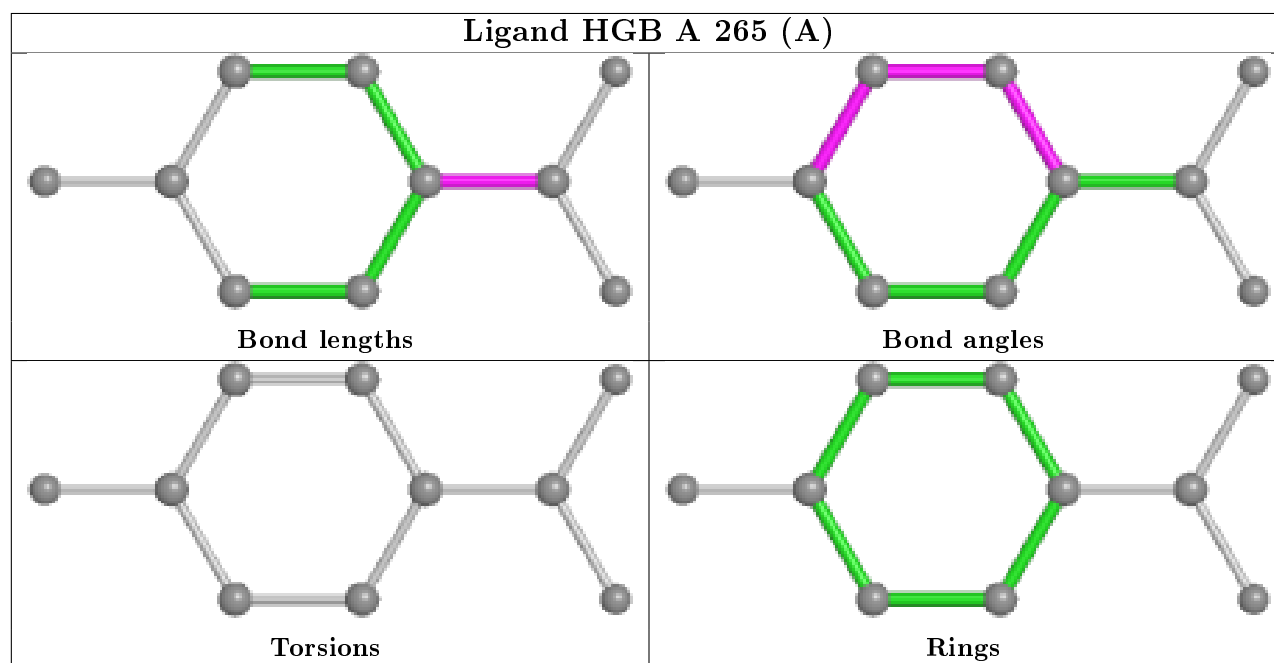
Mol	Chain	Res	Type	Atoms
5	A	1006	GOL	C1-C2-C3-O3
5	A	1004	GOL	O1-C1-C2-C3
5	A	1006	GOL	O2-C2-C3-O3
5	A	1004	GOL	O2-C2-C3-O3
4	A	1003	SUA	C11-C10-C9-C8
4	A	1003	SUA	C13-C10-C9-C8
4	A	1003	SUA	C4-C3-S1-O2
4	A	1003	SUA	C2-C3-S1-O2

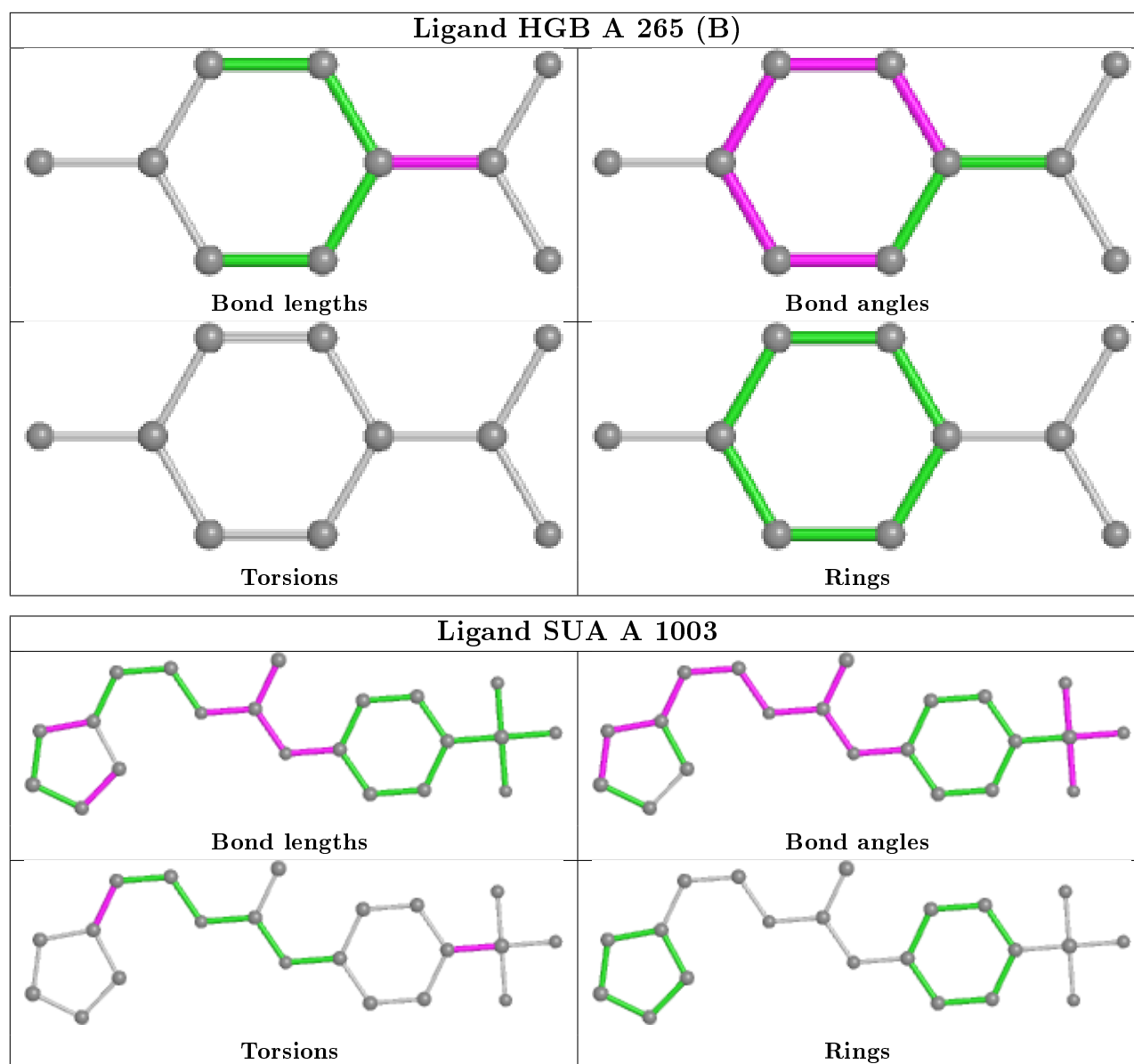
There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1006	GOL	2	0
5	A	1005	GOL	4	0
4	A	1003	SUA	5	0
5	A	1004	GOL	2	2

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	258/260 (99%)	-0.38	6 (2%) 60 43	6, 12, 26, 44	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	261	LYS	4.3
1	A	4	HIS	3.5
1	A	3	HIS	3.4
1	A	253	ASN	2.6
1	A	45	LYS	2.4
1	A	9	LYS	2.3

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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

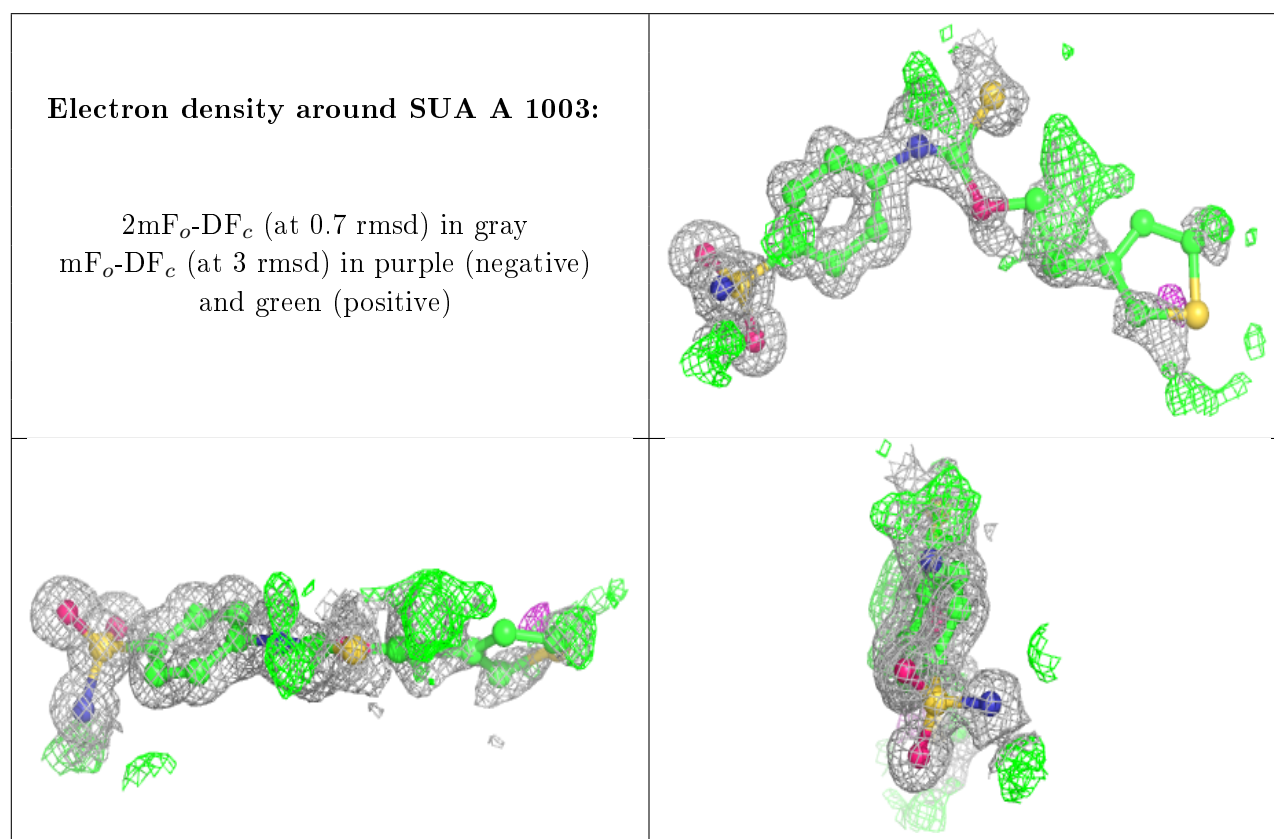
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	A	1005	6/6	0.85	0.18	21,26,28,28	0

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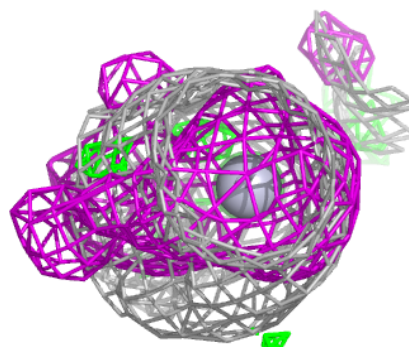
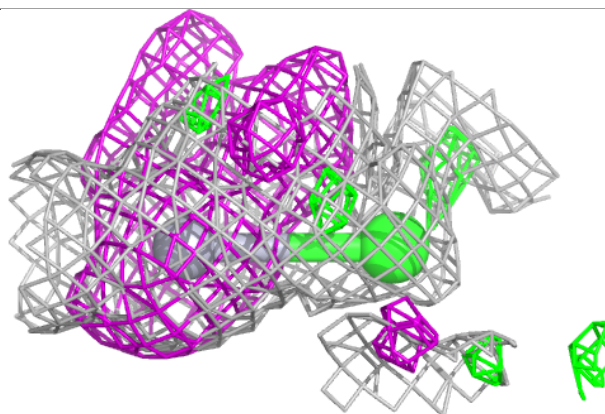
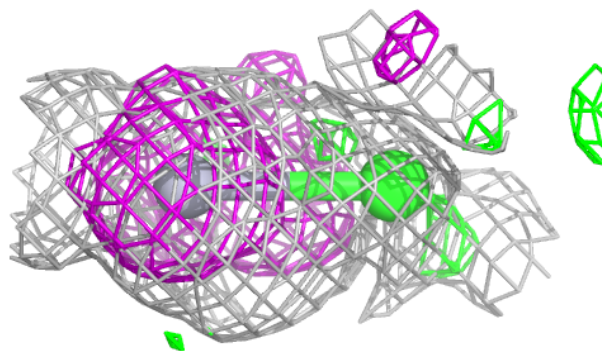
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	GOL	A	1004	6/6	0.90	0.16	21,24,43,57	0
5	GOL	A	1006	6/6	0.91	0.10	15,24,29,39	0
4	SUA	A	1003	21/21	0.99	0.07	6,14,40,59	10
2	HGB	A	265[B]	2/11	1.00	0.05	7,7,7,9	2
3	ZN	A	1002	1/1	1.00	0.03	5,5,5,5	0
2	HGB	A	265[A]	10/11	1.00	0.05	8,13,15,19	10

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.

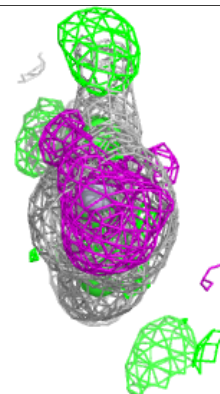
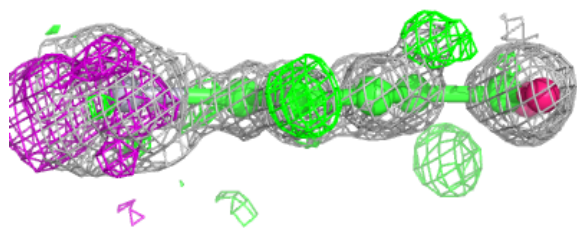
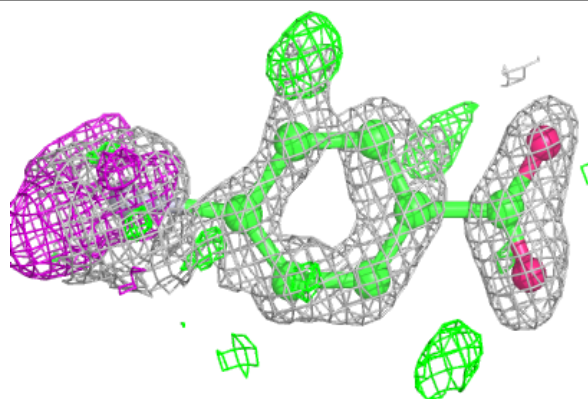


Electron density around HGB A 265 (B):

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

**Electron density around HGB A 265 (A):**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.