



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 08:19 PM UTC

PDB ID : 2R46 / pdb_00002r46
Title : Crystal structure of Escherichia coli Glycerol-3-phosphate Dehydrogenase in complex with 2-phosphopyruvic acid.
Authors : Yeh, J.I.; Du, S.; Chinte, U.
Deposited on : 2007-08-30
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 types of validation reports are described at

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

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

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

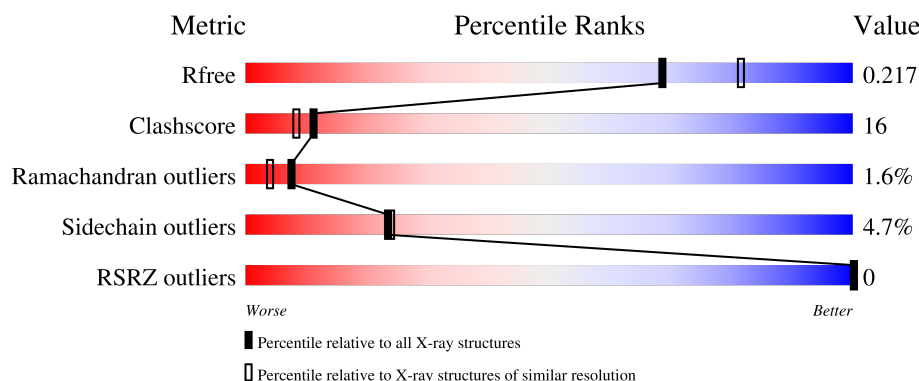
1 Overall quality at a glance

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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	501	 71% 24% 5% . .
1	B	501	 70% 24% 5% .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BOG	A	1949	X	-	-	-
2	BOG	A	800	X	-	X	-
6	PEP	A	700	-	X	X	-
6	PEP	B	700	-	X	-	-

2 Entry composition [i](#)

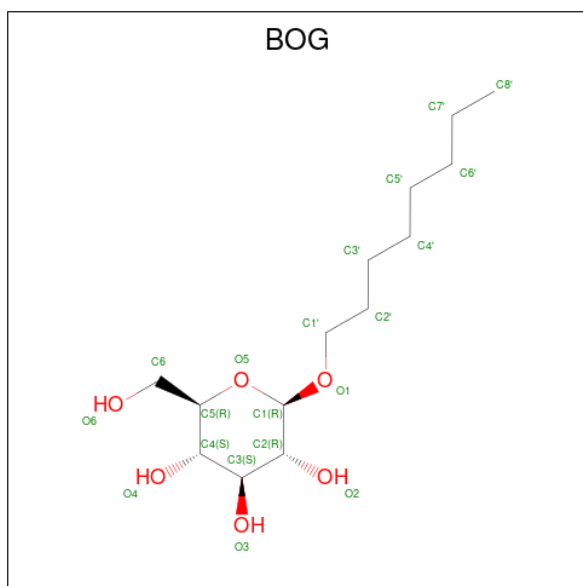
There are 8 unique types of molecules in this entry. The entry contains 8658 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aerobic glycerol-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	0	0
			3962	2515	705	729	13			
1	B	495	Total	C	N	O	S	0	0	0
			3962	2515	705	729	13			

- Molecule 2 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: $C_{14}H_{28}O_6$).



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

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



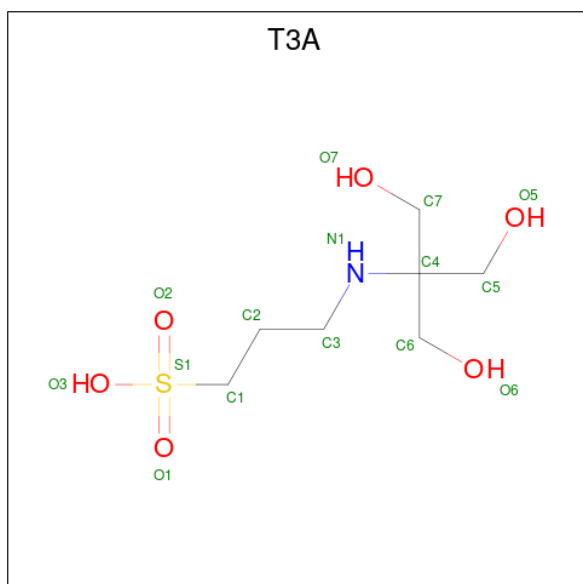
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		

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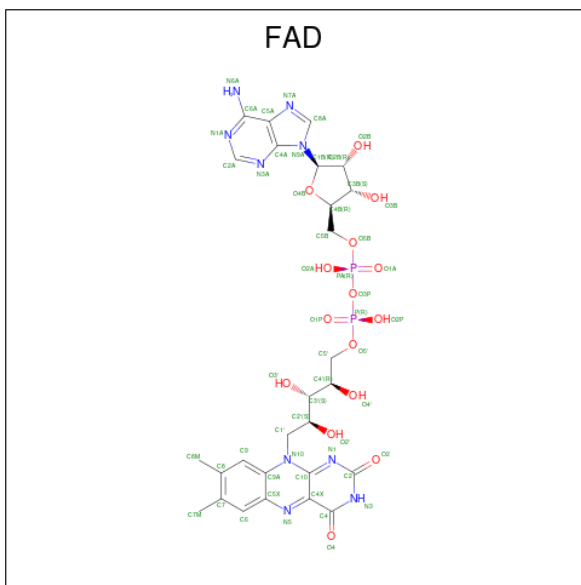
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total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

- Molecule 4 is N-(TRIS(HYDROXYMETHYL)METHYL)-3-AMINOPROPANESULFONIC ACID (CCD ID: T3A) (formula: $C_7H_{17}NO_6S$).



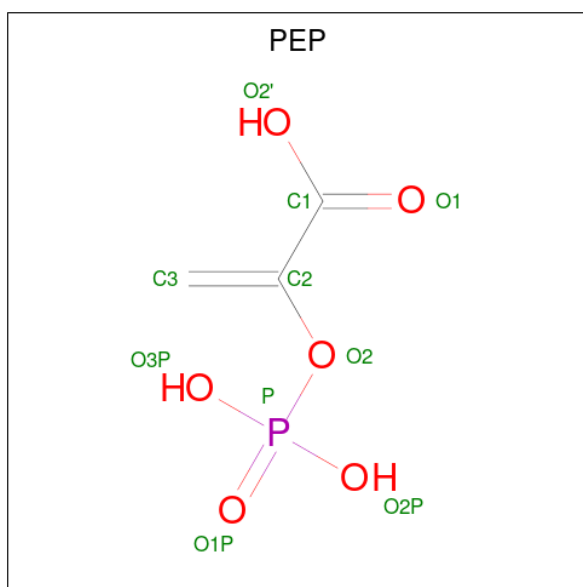
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 15	C 7	N 1	O 6	S 1	0	0
4	B	1	Total 15	C 7	N 1	O 6	S 1	0	0

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$).



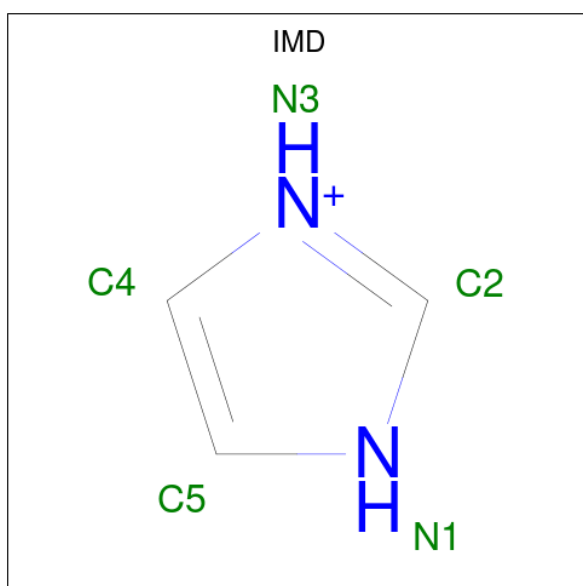
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
5	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 6 is PHOSPHOENOLPYRUVATE (CCD ID: PEP) (formula: $\text{C}_3\text{H}_5\text{O}_6\text{P}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	P	0	0
			10	3	6	1		
6	B	1	Total	C	O	P	0	0
			10	3	6	1		

- Molecule 7 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	N	0	0
			5	3	2		

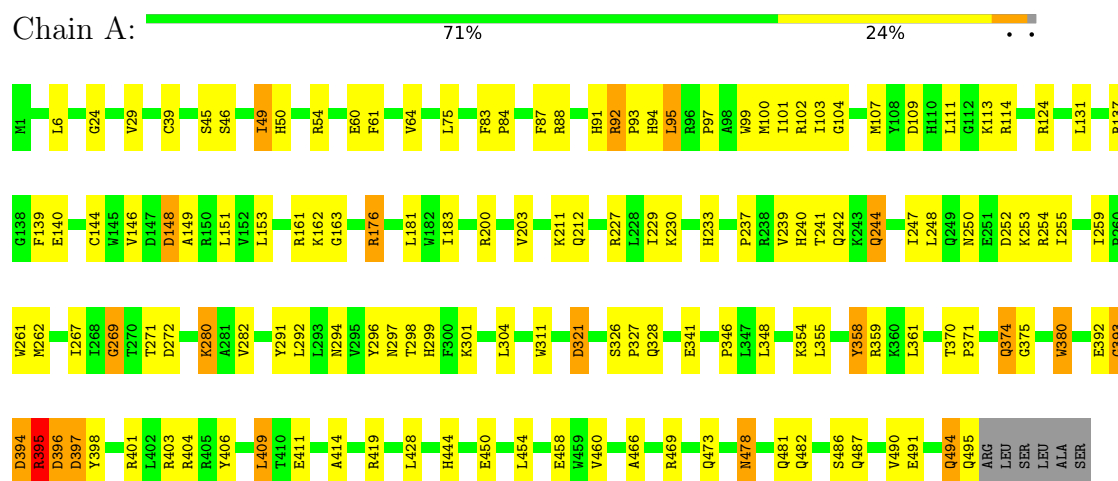
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	196	Total 196	O 196	0	0
8	B	189	Total 189	O 189	0	0

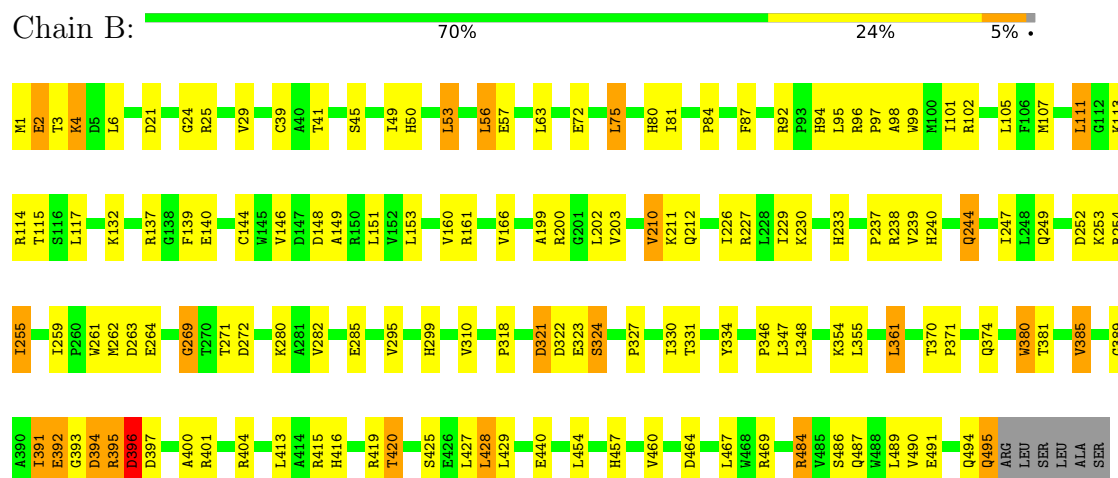
3 Residue-property plots

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: Aerobic glycerol-3-phosphate dehydrogenase



- Molecule 1: Aerobic glycerol-3-phosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	113.82Å 113.94Å 192.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.10 10.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (10.00-2.10) 99.7 (10.00-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.191 , 0.230 0.216 , 0.217	Depositor DCC
R_{free} test set	3706 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.480 for k,h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8658	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% 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: PEP, IMD, EDO, BOG, FAD, T3A

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	0.43	3/4057 (0.1%)	0.89	17/5494 (0.3%)
1	B	0.40	1/4057 (0.0%)	0.90	10/5494 (0.2%)
All	All	0.41	4/8114 (0.0%)	0.90	27/10988 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	291	TYR	C-N	7.20	1.43	1.33
1	B	269	GLY	C-O	-6.55	1.17	1.23
1	A	292	LEU	C-N	5.84	1.41	1.33
1	A	269	GLY	C-O	-5.07	1.18	1.23

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	395	ARG	N-CA-C	-12.95	95.76	111.69
1	A	395	ARG	N-CA-C	-11.20	98.03	112.93
1	A	259	ILE	N-CA-C	7.06	116.02	107.61
1	B	271	THR	CA-C-O	-6.87	113.82	121.51
1	B	132	LYS	CA-C-N	6.28	125.77	119.24
1	B	132	LYS	C-N-CA	6.28	125.77	119.24
1	B	259	ILE	N-CA-C	6.24	115.03	107.61
1	B	381	THR	N-CA-C	5.89	118.49	111.71
1	A	92	ARG	CA-C-N	5.82	127.11	119.84
1	A	92	ARG	C-N-CA	5.82	127.11	119.84
1	B	210	VAL	N-CA-C	5.63	116.27	110.36
1	A	398	TYR	N-CA-C	5.55	120.47	111.37
1	A	49	ILE	N-CA-C	-5.53	97.85	109.34
1	B	41	THR	N-CA-C	5.46	117.23	111.28
1	A	151	LEU	N-CA-C	-5.45	105.49	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	VAL	N-CA-C	5.27	116.90	108.89
1	A	271	THR	CA-C-O	-5.23	115.51	121.58
1	A	358	TYR	N-CA-C	5.23	116.98	111.28
1	B	151	LEU	N-CA-C	-5.21	105.68	111.36
1	A	291	TYR	CA-C-N	5.20	127.25	120.28
1	A	291	TYR	C-N-CA	5.20	127.25	120.28
1	A	237	PRO	N-CA-C	-5.18	103.06	111.14
1	B	29	VAL	N-CA-C	5.13	116.75	108.85
1	A	83	PHE	N-CA-C	5.11	112.82	108.22
1	A	61	PHE	N-CA-C	5.08	116.82	111.28
1	A	46	SER	N-CA-C	-5.02	107.12	113.20
1	A	148	ASP	N-CA-C	5.01	117.12	111.11

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	3962	0	3911	135	0
1	B	3962	0	3911	132	0
2	A	60	0	81	22	0
2	B	40	0	56	10	0
3	A	48	0	72	5	0
3	B	40	0	60	4	0
4	A	15	0	17	0	0
4	B	15	0	17	1	0
5	A	53	0	31	1	0
5	B	53	0	31	1	0
6	A	10	0	3	4	0
6	B	10	0	3	3	0
7	B	5	0	5	2	0
8	A	196	0	0	6	0
8	B	189	0	0	4	0
All	All	8658	0	8198	268	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 16.

All (268) 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:137:ARG:HH22	2:A:800:BOG:C5	1.62	1.10
1:A:137:ARG:HH22	2:A:800:BOG:H5	0.93	1.04
1:A:137:ARG:NH2	2:A:800:BOG:H5	1.77	0.99
1:A:254:ARG:HH11	6:A:700:PEP:H31	1.27	0.96
1:B:254:ARG:HH21	6:B:700:PEP:H31	1.32	0.94
1:B:105:LEU:HD13	2:B:900:BOG:O6	1.71	0.91
1:A:102:ARG:HD2	2:A:800:BOG:H3'1	1.53	0.90
1:A:254:ARG:NH1	6:A:700:PEP:H31	1.85	0.90
1:B:254:ARG:NH2	6:B:700:PEP:H31	1.86	0.90
1:A:297:ASN:HD21	1:A:304:LEU:H	1.20	0.89
1:B:391:ILE:HG13	1:B:392:GLU:H	1.38	0.87
1:B:395:ARG:HG2	1:B:419:ARG:HH12	1.42	0.85
1:B:415:ARG:HG2	1:B:419:ARG:HD2	1.58	0.84
1:A:102:ARG:HE	2:A:800:BOG:H1'1	1.43	0.83
1:A:395:ARG:NH1	1:A:419:ARG:HE	1.77	0.82
1:A:478:ASN:HD22	1:A:478:ASN:C	1.89	0.80
1:B:272:ASP:H	3:B:901:EDO:H21	1.45	0.80
1:A:395:ARG:NH1	1:A:395:ARG:HB2	1.96	0.80
1:A:478:ASN:ND2	1:A:481:GLN:H	1.79	0.80
2:A:1949:BOG:H8'3	2:A:1949:BOG:H2'1	1.63	0.80
1:B:324:SER:OG	1:B:330:ILE:HG12	1.80	0.80
1:A:272:ASP:H	3:A:1954:EDO:H12	1.48	0.79
1:A:254:ARG:HH11	6:A:700:PEP:C3	1.96	0.78
1:B:464:ASP:H	7:B:7068:IMD:HN3	1.28	0.78
1:B:105:LEU:CD1	2:B:900:BOG:O6	2.31	0.78
1:B:395:ARG:CG	1:B:419:ARG:HH12	1.98	0.77
1:A:103:ILE:HG22	1:A:107:MET:HE2	1.66	0.77
1:B:53:LEU:HD22	2:B:900:BOG:H5	1.65	0.76
1:B:395:ARG:HD3	1:B:419:ARG:HH22	1.49	0.76
1:B:1:MET:HE2	1:B:2:GLU:HG2	1.67	0.76
1:A:102:ARG:NH2	1:A:137:ARG:HH21	1.84	0.76
1:A:102:ARG:NE	2:A:800:BOG:H1'1	2.01	0.76
1:B:397:ASP:HA	1:B:400:ALA:HB3	1.68	0.76
1:B:416:HIS:O	1:B:420:THR:HG23	1.86	0.75
1:B:254:ARG:HH21	6:B:700:PEP:C3	2.00	0.74
1:A:93:PRO:HB3	2:A:1949:BOG:H5	1.69	0.74
1:B:96:ARG:HH11	1:B:249:GLN:HE21	1.36	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:393:GLY:CA	1:B:396:ASP:HB2	2.18	0.73
1:B:495:GLN:HE21	1:B:495:GLN:N	1.86	0.72
1:A:395:ARG:HB2	1:A:395:ARG:HH11	1.53	0.72
1:B:391:ILE:HG13	1:B:392:GLU:N	2.04	0.72
1:A:137:ARG:HH22	2:A:800:BOG:C6	2.03	0.72
1:B:322:ASP:OD2	1:B:331:THR:HG22	1.89	0.71
1:A:137:ARG:NH2	2:A:800:BOG:H62	2.06	0.71
1:B:393:GLY:HA3	1:B:396:ASP:HB2	1.73	0.71
1:B:6:LEU:HD11	1:B:203:VAL:HG23	1.74	0.69
1:B:374:GLN:CD	1:B:374:GLN:H	1.99	0.69
1:A:137:ARG:NH2	2:A:800:BOG:C6	2.56	0.69
1:B:395:ARG:O	1:B:397:ASP:N	2.26	0.68
1:A:102:ARG:NH2	1:A:137:ARG:NH2	2.41	0.68
1:A:137:ARG:NH2	2:A:800:BOG:C5	2.46	0.67
1:A:49:ILE:HB	1:A:144:CYS:HB2	1.77	0.66
1:A:397:ASP:HB3	1:A:401:ARG:HD3	1.78	0.66
1:B:49:ILE:HB	1:B:144:CYS:HB2	1.75	0.66
1:A:403:ARG:HD2	1:A:411:GLU:HA	1.79	0.65
1:A:54:ARG:HA	2:A:900:BOG:H61	1.79	0.65
1:B:102:ARG:HD2	2:B:800:BOG:H62	1.78	0.64
1:B:393:GLY:O	1:B:394:ASP:HB2	1.99	0.63
1:B:202:LEU:HD13	1:B:203:VAL:N	2.14	0.62
1:A:94:HIS:CE1	1:A:95:LEU:HD13	2.35	0.62
1:B:72:GLU:OE1	1:B:115:THR:HG22	1.99	0.62
1:B:50:HIS:CE1	1:B:354:LYS:HZ3	2.17	0.62
1:A:107:MET:O	1:A:111:LEU:HB2	2.00	0.62
1:B:395:ARG:HG2	1:B:419:ARG:NH1	2.15	0.61
1:B:393:GLY:O	1:B:394:ASP:CB	2.47	0.61
1:A:478:ASN:HD21	1:A:481:GLN:H	1.47	0.61
1:A:294:ASN:O	1:A:298:THR:HG23	2.00	0.61
1:B:244:GLN:NE2	1:B:244:GLN:H	1.99	0.61
1:B:1:MET:HE3	1:B:2:GLU:H	1.66	0.61
1:A:93:PRO:CB	2:A:1949:BOG:H5	2.30	0.61
1:B:57:GLU:HB2	2:B:900:BOG:O2	2.01	0.61
1:A:102:ARG:CD	2:A:800:BOG:H1'1	2.31	0.60
1:A:428:LEU:C	1:A:428:LEU:HD23	2.27	0.60
1:A:49:ILE:HD11	1:A:146:VAL:HB	1.82	0.60
1:A:393:GLY:O	1:A:394:ASP:CB	2.49	0.60
1:B:334:TYR:CD2	1:B:361:LEU:HD22	2.36	0.60
1:A:6:LEU:HD11	1:A:203:VAL:HG23	1.83	0.60
1:A:490:VAL:O	1:A:494:GLN:HG2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:395:ARG:CD	1:B:419:ARG:HH22	2.15	0.60
1:B:428:LEU:HD12	1:B:428:LEU:O	2.01	0.60
1:B:392:GLU:HG2	1:B:401:ARG:HH12	1.68	0.59
1:B:238:ARG:HG3	1:B:264:GLU:C	2.27	0.58
1:A:487:GLN:O	1:A:491:GLU:HG3	2.03	0.58
1:B:50:HIS:CE1	1:B:354:LYS:NZ	2.71	0.58
1:A:24:GLY:HA2	1:A:162:LYS:HD3	1.85	0.58
1:A:395:ARG:CZ	1:A:419:ARG:HE	2.17	0.58
1:B:24:GLY:HA3	1:B:380:TRP:CZ2	2.39	0.58
1:B:428:LEU:HD12	1:B:428:LEU:C	2.29	0.57
1:A:241:THR:HG22	1:A:241:THR:O	2.03	0.57
1:A:102:ARG:HD2	2:A:800:BOG:C1'	2.34	0.57
1:B:487:GLN:O	1:B:491:GLU:HG3	2.04	0.57
1:A:478:ASN:C	1:A:478:ASN:ND2	2.62	0.57
1:A:54:ARG:HA	2:A:900:BOG:C6	2.35	0.56
1:B:397:ASP:O	1:B:401:ARG:N	2.23	0.56
1:B:229:ILE:HD12	1:B:327:PRO:HB3	1.88	0.56
1:B:324:SER:CB	1:B:330:ILE:HG12	2.35	0.56
1:A:102:ARG:HD2	2:A:800:BOG:H1'1	1.87	0.56
1:B:247:ILE:HG13	1:B:255:ILE:HG23	1.87	0.55
1:A:200:ARG:O	1:A:346:PRO:HD2	2.07	0.55
1:A:444:HIS:HE1	8:A:7156:HOH:O	1.88	0.55
1:A:102:ARG:HH22	1:A:137:ARG:HH21	1.52	0.55
1:A:374:GLN:CG	1:A:375:GLY:H	2.19	0.55
1:B:113:LYS:HB2	8:B:7227:HOH:O	2.05	0.55
1:A:403:ARG:HD3	8:A:7254:HOH:O	2.08	0.54
1:A:406:TYR:HB2	1:A:409:LEU:HD22	1.90	0.54
1:B:107:MET:O	1:B:111:LEU:HB2	2.07	0.53
1:B:113:LYS:HB2	1:B:113:LYS:NZ	2.24	0.53
1:B:252:ASP:O	1:B:253:LYS:HB2	2.08	0.53
1:A:102:ARG:HD2	2:A:800:BOG:C3'	2.34	0.53
1:A:478:ASN:ND2	1:A:481:GLN:HG3	2.24	0.53
1:B:99:TRP:HB3	2:B:800:BOG:H1'2	1.91	0.52
1:A:428:LEU:HD11	1:A:450:GLU:HG3	1.91	0.52
3:A:1956:EDO:O1	1:B:161:ARG:HB2	2.09	0.52
1:B:153:LEU:HD13	1:B:153:LEU:C	2.35	0.52
1:A:478:ASN:HD21	1:A:481:GLN:HG3	1.74	0.51
1:B:230:LYS:NZ	1:B:285:GLU:HG3	2.25	0.51
1:B:272:ASP:N	3:B:901:EDO:H21	2.21	0.51
1:B:440:GLU:OE1	1:B:484:ARG:HD3	2.11	0.51
1:A:88:ARG:NE	1:A:244:GLN:HG3	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:TRP:HA	2:A:800:BOG:H1'2	1.92	0.51
1:B:391:ILE:O	1:B:392:GLU:CB	2.59	0.51
1:B:49:ILE:HD11	1:B:146:VAL:HB	1.91	0.51
1:A:24:GLY:HA3	1:A:380:TRP:CZ2	2.46	0.51
1:A:240:HIS:CE1	1:A:242:GLN:HB2	2.46	0.51
1:A:137:ARG:NH1	2:A:1949:BOG:O2	2.44	0.50
1:A:252:ASP:O	1:A:253:LYS:HB2	2.10	0.50
1:A:227:ARG:HG2	1:A:321:ASP:HB2	1.94	0.50
1:A:247:ILE:HG23	1:A:247:ILE:O	2.12	0.50
1:B:226:ILE:HD11	1:B:318:PRO:C	2.37	0.50
1:A:444:HIS:HD2	1:A:473:GLN:OE1	1.95	0.50
1:B:200:ARG:O	1:B:346:PRO:HD2	2.12	0.50
1:A:104:GLY:HA2	1:A:107:MET:HE3	1.93	0.49
1:A:403:ARG:NH1	1:A:409:LEU:O	2.46	0.49
1:A:454:LEU:HB3	1:A:460:VAL:HG21	1.94	0.49
1:A:494:GLN:HB3	1:A:495:GLN:OE1	2.12	0.49
1:B:202:LEU:HD13	1:B:202:LEU:C	2.38	0.49
1:B:92:ARG:HD2	1:B:299:HIS:HE1	1.77	0.49
1:B:370:THR:HB	1:B:371:PRO:HD3	1.94	0.49
1:B:98:ALA:HB3	2:B:800:BOG:H3	1.94	0.49
1:A:326:SER:OG	1:A:328:GLN:HG2	2.13	0.48
1:A:361:LEU:C	1:A:361:LEU:HD23	2.39	0.48
1:A:176:ARG:HD2	1:A:183:ILE:HB	1.94	0.48
1:A:229:ILE:HD12	1:A:327:PRO:HB3	1.94	0.48
1:A:392:GLU:HB2	1:A:401:ARG:NH1	2.28	0.48
1:B:21:ASP:O	1:B:25:ARG:HG3	2.12	0.48
1:B:212:GLN:NE2	8:B:7127:HOH:O	2.46	0.48
1:A:395:ARG:HH22	1:A:419:ARG:HG3	1.78	0.48
1:B:1:MET:CE	1:B:2:GLU:HG2	2.39	0.48
1:B:391:ILE:CG1	1:B:392:GLU:H	2.09	0.48
1:B:396:ASP:OD2	1:B:396:ASP:N	2.47	0.48
1:A:102:ARG:HH12	1:A:139:PHE:HE1	1.61	0.48
1:B:94:HIS:CE1	1:B:95:LEU:HG	2.48	0.48
1:A:148:ASP:OD1	1:A:149:ALA:N	2.46	0.48
1:A:255:ILE:HD12	1:A:255:ILE:C	2.38	0.48
8:A:7174:HOH:O	1:B:457:HIS:HE1	1.95	0.48
1:B:101:ILE:HA	2:B:900:BOG:H61	1.95	0.47
1:B:454:LEU:HB3	1:B:460:VAL:HG21	1.94	0.47
1:A:395:ARG:HH12	1:A:419:ARG:HE	1.58	0.47
1:B:202:LEU:HD12	1:B:347:LEU:CD1	2.44	0.47
1:B:361:LEU:HD12	1:B:361:LEU:C	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:VAL:O	1:B:299:HIS:HD2	1.97	0.47
1:A:348:LEU:HD23	1:A:348:LEU:C	2.40	0.47
1:B:92:ARG:HD2	1:B:299:HIS:CE1	2.49	0.47
1:A:397:ASP:HB3	1:A:401:ARG:HH11	1.79	0.47
1:A:374:GLN:CG	1:A:375:GLY:N	2.78	0.46
1:A:97:PRO:HG2	1:A:100:MET:HB2	1.97	0.46
1:B:72:GLU:OE2	1:B:114:ARG:HA	2.15	0.46
1:A:75:LEU:HD21	1:A:84:PRO:HG3	1.98	0.46
1:A:94:HIS:NE2	1:A:95:LEU:HD13	2.31	0.46
1:A:102:ARG:HG2	1:A:102:ARG:HH11	1.81	0.46
1:A:255:ILE:HD12	1:A:255:ILE:O	2.16	0.46
1:A:181:LEU:HD23	8:A:7239:HOH:O	2.16	0.46
1:A:272:ASP:H	3:A:1954:EDO:C1	2.23	0.46
1:A:60:GLU:O	1:A:64:VAL:HG23	2.15	0.46
1:B:227:ARG:HG2	1:B:321:ASP:CB	2.45	0.46
1:A:478:ASN:O	1:A:482:GLN:HG3	2.16	0.46
1:B:101:ILE:HG21	1:B:139:PHE:CD2	2.51	0.46
1:B:427:LEU:HD12	8:B:7217:HOH:O	2.16	0.46
1:B:102:ARG:NH2	1:B:137:ARG:HH12	2.13	0.46
1:B:233:HIS:CD2	1:B:269:GLY:HA3	2.51	0.46
1:A:153:LEU:HD21	1:A:458:GLU:CD	2.40	0.45
1:B:210:VAL:HG13	1:B:211:LYS:N	2.31	0.45
1:B:391:ILE:O	1:B:392:GLU:HB3	2.15	0.45
1:A:176:ARG:HH11	1:A:176:ARG:HG3	1.82	0.45
1:B:97:PRO:HB2	1:B:99:TRP:CD1	2.51	0.45
1:B:395:ARG:HH11	1:B:395:ARG:HG3	1.81	0.45
1:A:241:THR:O	1:A:241:THR:CG2	2.64	0.45
1:A:406:TYR:O	1:A:409:LEU:HB2	2.16	0.45
1:B:39:CYS:HA	1:B:469:ARG:HD3	1.97	0.45
1:A:248:LEU:O	1:A:255:ILE:HA	2.17	0.45
1:A:50:HIS:NE2	1:A:354:LYS:NZ	2.47	0.45
1:A:102:ARG:NH1	1:A:139:PHE:HE1	2.14	0.45
1:A:161:ARG:HH21	3:A:1953:EDO:H11	1.80	0.45
1:B:87:PHE:O	1:B:140:GLU:HA	2.17	0.45
1:B:227:ARG:HG2	1:B:321:ASP:HB2	1.98	0.45
1:B:361:LEU:C	1:B:361:LEU:CD1	2.90	0.45
1:A:403:ARG:HG3	1:A:414:ALA:HB2	1.98	0.45
1:B:96:ARG:HD2	1:B:249:GLN:NE2	2.32	0.45
1:A:87:PHE:O	1:A:140:GLU:HA	2.16	0.44
1:B:490:VAL:HG13	1:B:494:GLN:NE2	2.32	0.44
1:A:153:LEU:CD2	1:A:458:GLU:CD	2.91	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:SER:HB2	5:B:600:FAD:C4X	2.47	0.44
1:B:419:ARG:HG3	1:B:419:ARG:HH11	1.82	0.44
1:B:419:ARG:HH11	1:B:419:ARG:CG	2.30	0.44
1:A:45:SER:HB2	5:A:600:FAD:C4X	2.47	0.44
1:A:163:GLY:HA3	1:B:404:ARG:HG2	1.98	0.44
1:A:233:HIS:CD2	1:A:269:GLY:HA3	2.52	0.44
1:B:148:ASP:OD1	1:B:149:ALA:N	2.50	0.44
1:B:50:HIS:HE1	1:B:354:LYS:NZ	2.16	0.44
1:A:91:HIS:CD2	1:A:93:PRO:HD3	2.52	0.44
1:A:162:LYS:HE3	1:A:380:TRP:CH2	2.53	0.44
1:B:495:GLN:N	1:B:495:GLN:NE2	2.60	0.44
1:A:267:ILE:HD12	8:A:7242:HOH:O	2.18	0.44
1:A:261:TRP:CE2	1:A:262:MET:HE2	2.53	0.44
1:B:72:GLU:OE2	1:B:117:LEU:HD23	2.18	0.44
1:A:261:TRP:CZ2	1:A:262:MET:HE2	2.53	0.43
1:B:202:LEU:HD12	1:B:347:LEU:HD13	2.00	0.43
1:A:101:ILE:HG21	1:A:139:PHE:CD2	2.54	0.43
1:A:102:ARG:CD	2:A:800:BOG:C1'	2.95	0.43
1:B:107:MET:HB3	1:B:111:LEU:HD22	2.00	0.43
1:B:374:GLN:CD	1:B:374:GLN:N	2.71	0.43
1:B:486:SER:O	1:B:490:VAL:HG23	2.17	0.43
1:A:49:ILE:CD1	1:A:146:VAL:HB	2.48	0.43
1:A:230:LYS:HD3	1:A:282:VAL:HG23	2.01	0.43
1:A:280:LYS:HA	1:A:280:LYS:HE3	2.01	0.43
1:B:230:LYS:HD3	1:B:282:VAL:HG23	2.00	0.43
1:B:80:HIS:CD2	1:B:81:ILE:HG23	2.53	0.43
1:A:254:ARG:NH1	6:A:700:PEP:C3	2.66	0.43
1:B:457:HIS:HA	4:B:7066:T3A:H22	2.01	0.43
1:A:244:GLN:H	1:A:244:GLN:CD	2.26	0.42
1:A:131:LEU:O	1:A:301:LYS:HE3	2.19	0.42
1:B:322:ASP:OD1	1:B:330:ILE:HG23	2.19	0.42
1:B:487:GLN:HE21	1:B:491:GLU:CG	2.33	0.42
1:A:460:VAL:HG11	1:A:466:ALA:HB2	2.00	0.42
1:B:385:VAL:O	1:B:385:VAL:CG2	2.66	0.42
1:B:247:ILE:HG13	1:B:255:ILE:CG2	2.48	0.42
1:A:109:ASP:HB3	1:A:114:ARG:HH21	1.84	0.42
1:B:75:LEU:HD21	1:B:84:PRO:HG3	2.02	0.42
1:A:39:CYS:HA	1:A:469:ARG:HD3	2.02	0.42
1:A:370:THR:HB	1:A:371:PRO:HD3	2.00	0.42
1:B:56:LEU:HB3	2:B:900:BOG:H1'2	2.01	0.42
1:B:395:ARG:CG	1:B:419:ARG:HH22	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:VAL:HG23	1:A:240:HIS:CD2	2.55	0.41
1:B:53:LEU:CD2	2:B:900:BOG:H5	2.42	0.41
1:A:211:LYS:HD3	8:A:7218:HOH:O	2.21	0.41
1:B:237:PRO:O	1:B:238:ARG:C	2.64	0.41
1:B:484:ARG:HH21	1:B:487:GLN:HG3	1.85	0.41
1:A:97:PRO:HB2	1:A:99:TRP:CD1	2.54	0.41
1:A:374:GLN:HG2	1:A:375:GLY:N	2.36	0.41
1:B:425:SER:O	1:B:429:LEU:HD23	2.20	0.41
1:A:311:TRP:HA	3:A:7069:EDO:H21	2.03	0.41
1:B:239:VAL:HG23	1:B:240:HIS:CD2	2.56	0.41
1:A:358:TYR:CG	1:A:359:ARG:N	2.88	0.41
1:A:393:GLY:O	1:A:394:ASP:HB2	2.19	0.41
1:A:296:TYR:CD1	1:A:296:TYR:C	2.99	0.41
1:A:393:GLY:O	1:A:394:ASP:HB3	2.21	0.41
1:A:486:SER:O	1:A:490:VAL:HG23	2.21	0.41
1:B:160:VAL:HG22	3:B:7067:EDO:H21	2.03	0.41
1:B:348:LEU:C	1:B:348:LEU:HD23	2.45	0.41
1:B:4:LYS:O	1:B:199:ALA:HA	2.21	0.41
7:B:7068:IMD:H4	8:B:7195:HOH:O	2.19	0.41
1:A:92:ARG:HD2	1:A:299:HIS:NE2	2.36	0.40
1:B:166:VAL:O	3:B:7067:EDO:H11	2.21	0.40
1:B:1:MET:O	1:B:2:GLU:C	2.64	0.40
1:B:261:TRP:CZ2	1:B:262:MET:HE2	2.56	0.40
1:B:392:GLU:HG2	1:B:401:ARG:NH1	2.33	0.40
1:A:92:ARG:HA	1:A:93:PRO:HD2	1.84	0.40
1:B:322:ASP:O	1:B:323:GLU:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	493/501 (98%)	466 (94%)	21 (4%)	6 (1%)	10	7
1	B	493/501 (98%)	465 (94%)	18 (4%)	10 (2%)	6	2
All	All	986/1002 (98%)	931 (94%)	39 (4%)	16 (2%)	7	4

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	GLN
1	A	394	ASP
1	B	392	GLU
1	B	394	ASP
1	B	396	ASP
1	B	2	GLU
1	A	380	TRP
1	A	393	GLY
1	A	396	ASP
1	B	321	ASP
1	B	3	THR
1	B	380	TRP
1	A	321	ASP
1	B	263	ASP
1	B	391	ILE
1	B	389	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	414/419 (99%)	397 (96%)	17 (4%)	27	29
1	B	414/419 (99%)	392 (95%)	22 (5%)	20	19
All	All	828/838 (99%)	789 (95%)	39 (5%)	23	24

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	LEU
1	A	113	LYS
1	A	124	ARG
1	A	176	ARG
1	A	212	GLN
1	A	244	GLN
1	A	250	ASN
1	A	280	LYS
1	A	341	GLU
1	A	355	LEU
1	A	395	ARG
1	A	396	ASP
1	A	397	ASP
1	A	404	ARG
1	A	409	LEU
1	A	478	ASN
1	A	494	GLN
1	B	4	LYS
1	B	53	LEU
1	B	56	LEU
1	B	63	LEU
1	B	75	LEU
1	B	111	LEU
1	B	244	GLN
1	B	255	ILE
1	B	280	LYS
1	B	310	VAL
1	B	324	SER
1	B	355	LEU
1	B	361	LEU
1	B	385	VAL
1	B	396	ASP
1	B	413	LEU
1	B	420	THR
1	B	428	LEU
1	B	467	LEU
1	B	484	ARG
1	B	489	LEU
1	B	495	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	35	GLN
1	A	128	ASN
1	A	179	ASN
1	A	212	GLN
1	A	242	GLN
1	A	250	ASN
1	A	297	ASN
1	A	424	ASN
1	A	444	HIS
1	A	473	GLN
1	A	478	ASN
1	A	482	GLN
1	A	494	GLN
1	B	50	HIS
1	B	91	HIS
1	B	94	HIS
1	B	198	GLN
1	B	212	GLN
1	B	219	HIS
1	B	244	GLN
1	B	249	GLN
1	B	299	HIS
1	B	364	HIS
1	B	424	ASN
1	B	457	HIS
1	B	473	GLN
1	B	487	GLN
1	B	494	GLN
1	B	495	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

34 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$
3	EDO	A	1953	-	3,3,3	0.47	0	2,2,2	0.35	0
3	EDO	B	903	-	3,3,3	0.41	0	2,2,2	0.37	0
3	EDO	B	904	-	3,3,3	0.48	0	2,2,2	0.33	0
3	EDO	A	1952	-	3,3,3	0.41	0	2,2,2	0.37	0
7	IMD	B	7068	-	5,5,5	0.35	0	5,5,5	0.38	0
3	EDO	A	7067	-	3,3,3	0.42	0	2,2,2	0.40	0
3	EDO	A	7070	-	3,3,3	0.33	0	2,2,2	0.65	0
3	EDO	A	1951	-	3,3,3	0.50	0	2,2,2	0.33	0
3	EDO	A	7069	-	3,3,3	0.47	0	2,2,2	0.35	0
3	EDO	B	902	-	3,3,3	0.44	0	2,2,2	0.35	0
3	EDO	B	908	-	3,3,3	0.39	0	2,2,2	0.39	0
6	PEP	A	700	-	9,9,9	5.24	4 (44%)	11,13,13	6.47	5 (45%)
2	BOG	B	900	-	20,20,20	0.46	0	25,25,25	0.58	0
2	BOG	A	800	-	20,20,20	3.41	3 (15%)	25,25,25	4.40	8 (32%)
2	BOG	A	900	-	20,20,20	0.47	0	25,25,25	0.59	0
3	EDO	A	1956	-	3,3,3	0.44	0	2,2,2	0.35	0
3	EDO	B	7067	-	3,3,3	0.43	0	2,2,2	0.37	0
2	BOG	A	1949	-	20,20,20	0.54	0	25,25,25	2.49	6 (24%)
3	EDO	B	905	-	3,3,3	0.47	0	2,2,2	0.33	0
6	PEP	B	700	-	9,9,9	5.53	5 (55%)	11,13,13	5.60	6 (54%)
3	EDO	A	7068	-	3,3,3	0.46	0	2,2,2	0.35	0
5	FAD	B	600	-	58,58,58	1.45	5 (8%)	85,89,89	1.68	17 (20%)
4	T3A	B	7066	-	14,14,14	1.12	1 (7%)	19,19,19	1.68	6 (31%)
3	EDO	A	1957	-	3,3,3	0.43	0	2,2,2	0.37	0
5	FAD	A	600	-	58,58,58	1.50	11 (18%)	85,89,89	1.72	18 (21%)
3	EDO	B	906	-	3,3,3	0.46	0	2,2,2	0.36	0
2	BOG	B	800	-	20,20,20	0.60	1 (5%)	25,25,25	2.39	10 (40%)
3	EDO	A	1955	-	3,3,3	0.41	0	2,2,2	0.37	0
3	EDO	B	901	-	3,3,3	0.44	0	2,2,2	0.36	0
4	T3A	A	7066	-	14,14,14	1.10	1 (7%)	19,19,19	1.68	6 (31%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	A	1950	-	3,3,3	0.46	0	2,2,2	0.34	0
3	EDO	A	1954	-	3,3,3	0.43	0	2,2,2	0.36	0
3	EDO	B	7069	-	3,3,3	0.45	0	2,2,2	0.88	0
3	EDO	B	907	-	3,3,3	0.43	0	2,2,2	0.35	0

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
3	EDO	A	1953	-	-	1/1/1/1	-
3	EDO	B	903	-	-	1/1/1/1	-
3	EDO	B	904	-	-	1/1/1/1	-
3	EDO	A	1952	-	-	1/1/1/1	-
7	IMD	B	7068	-	-	-	0/1/1/1
3	EDO	A	7067	-	-	0/1/1/1	-
3	EDO	A	7070	-	-	1/1/1/1	-
3	EDO	A	1951	-	-	1/1/1/1	-
3	EDO	A	7069	-	-	1/1/1/1	-
3	EDO	B	902	-	-	1/1/1/1	-
3	EDO	B	908	-	-	0/1/1/1	-
6	PEP	A	700	-	-	4/9/9/9	-
2	BOG	B	900	-	-	3/11/31/31	0/1/1/1
2	BOG	A	800	-	2/2/5/5	3/11/31/31	0/1/1/1
2	BOG	A	900	-	-	6/11/31/31	0/1/1/1
3	EDO	A	1956	-	-	1/1/1/1	-
3	EDO	B	7067	-	-	1/1/1/1	-
2	BOG	A	1949	-	2/2/5/5	5/11/31/31	0/1/1/1
3	EDO	B	905	-	-	0/1/1/1	-
6	PEP	B	700	-	-	4/9/9/9	-
3	EDO	A	7068	-	-	0/1/1/1	-
5	FAD	B	600	-	-	1/34/50/50	0/6/6/6
4	T3A	B	7066	-	-	0/18/18/18	-
3	EDO	A	1957	-	-	0/1/1/1	-
5	FAD	A	600	-	-	2/34/50/50	0/6/6/6
3	EDO	B	906	-	-	1/1/1/1	-
2	BOG	B	800	-	-	6/11/31/31	0/1/1/1
3	EDO	A	1955	-	-	1/1/1/1	-
3	EDO	B	901	-	-	1/1/1/1	-
4	T3A	A	7066	-	-	0/18/18/18	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1950	-	-	1/1/1/1	-
3	EDO	A	1954	-	-	1/1/1/1	-
3	EDO	B	7069	-	-	1/1/1/1	-
3	EDO	B	907	-	-	1/1/1/1	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	700	PEP	C2-C1	12.08	1.61	1.49
2	A	800	BOG	C4-C3	-11.35	1.22	1.52
6	A	700	PEP	C2-C1	11.26	1.60	1.49
6	B	700	PEP	C3-C2	9.75	1.57	1.31
6	A	700	PEP	C3-C2	9.28	1.56	1.31
2	A	800	BOG	C4-C5	-8.39	1.35	1.53
5	B	600	FAD	C4X-N5	5.37	1.42	1.30
2	A	800	BOG	O4-C4	5.10	1.55	1.43
5	A	600	FAD	C4X-N5	4.17	1.39	1.30
5	B	600	FAD	C5A-N7A	-4.05	1.31	1.39
5	B	600	FAD	C10-N1	3.75	1.40	1.33
6	B	700	PEP	O2-C2	3.64	1.49	1.39
5	A	600	FAD	C10-N1	3.64	1.40	1.33
6	A	700	PEP	O2-C2	3.60	1.49	1.39
6	A	700	PEP	O2'-C1	3.56	1.40	1.30
5	A	600	FAD	C2A-N1A	3.12	1.39	1.33
5	B	600	FAD	C6-C7	-2.96	1.35	1.39
6	B	700	PEP	O2'-C1	2.96	1.38	1.30
5	A	600	FAD	C4A-N3A	2.91	1.39	1.34
5	B	600	FAD	C2A-N3A	2.90	1.39	1.33
5	A	600	FAD	C5'-C4'	2.83	1.55	1.51
5	A	600	FAD	C2A-N3A	2.80	1.38	1.33
5	A	600	FAD	P-O3P	2.71	1.62	1.59
5	A	600	FAD	O4-C4	2.59	1.28	1.23
5	A	600	FAD	C7M-C7	-2.55	1.46	1.51
6	B	700	PEP	P-O3P	-2.46	1.45	1.54
5	A	600	FAD	C1'-N10	2.42	1.53	1.47
4	A	7066	T3A	C1-S1	2.38	1.80	1.77
4	B	7066	T3A	C1-S1	2.29	1.80	1.77
5	A	600	FAD	C5A-N7A	-2.13	1.35	1.39
2	B	800	BOG	O1-C1	2.02	1.43	1.40

All (82) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	BOG	C3-C4-C5	17.26	141.52	110.23
6	A	700	PEP	O1-C1-C2	-13.22	101.85	121.79
6	A	700	PEP	O2'-C1-C2	12.84	135.81	113.91
6	B	700	PEP	O2'-C1-C2	11.64	133.77	113.91
6	B	700	PEP	O1-C1-C2	-10.34	106.20	121.79
6	A	700	PEP	O2-C2-C3	-10.06	105.62	124.85
2	A	1949	BOG	O2-C2-C1	9.74	133.29	110.08
6	B	700	PEP	O2-C2-C3	-9.08	107.50	124.85
5	A	600	FAD	N3A-C2A-N1A	-7.46	117.29	128.58
2	A	800	BOG	O5-C5-C4	-7.35	96.46	109.70
2	A	800	BOG	O3-C3-C4	7.08	127.07	110.38
2	A	800	BOG	C6-C5-C4	-6.33	97.47	113.02
5	B	600	FAD	N3A-C2A-N1A	-6.26	119.10	128.58
2	B	800	BOG	O1-C1-C2	5.04	115.93	108.27
2	B	800	BOG	O5-C1-C2	-4.56	101.00	110.37
5	A	600	FAD	N9A-C8A-N7A	-4.30	107.84	113.94
5	B	600	FAD	C5A-N7A-C8A	4.29	110.19	103.45
2	A	1949	BOG	O3-C3-C4	4.29	120.49	110.38
5	B	600	FAD	C5A-C4A-N3A	-4.28	120.83	126.72
2	B	800	BOG	C1'-O1-C1	-4.16	106.57	113.68
5	B	600	FAD	N9A-C8A-N7A	-4.07	108.16	113.94
2	B	800	BOG	O5-C5-C4	-3.85	102.77	109.70
5	A	600	FAD	C5A-N7A-C8A	3.71	109.28	103.45
5	B	600	FAD	C2A-N3A-C4A	3.63	120.69	111.83
5	A	600	FAD	C2A-N3A-C4A	3.53	120.44	111.83
5	A	600	FAD	C5A-C4A-N3A	-3.48	121.93	126.72
5	A	600	FAD	C10-C4X-N5	-3.47	117.72	124.81
4	A	7066	T3A	C2-C1-S1	-3.46	107.94	113.25
2	A	800	BOG	O5-C1-O1	3.44	118.16	110.04
4	B	7066	T3A	C2-C1-S1	-3.36	108.10	113.25
5	B	600	FAD	C4A-C5A-N7A	-3.22	106.90	110.58
2	A	800	BOG	C1'-O1-C1	3.16	119.07	113.68
2	A	800	BOG	O5-C5-C6	3.06	114.03	106.44
2	A	1949	BOG	O5-C5-C6	3.03	113.94	106.44
5	A	600	FAD	N3A-C4A-N9A	3.00	132.27	127.17
4	A	7066	T3A	C3-N1-C4	-2.97	111.81	116.17
2	B	800	BOG	O5-C1-O1	-2.96	103.06	110.04
5	B	600	FAD	C5X-C9A-N10	2.93	120.62	117.97
2	B	800	BOG	O2-C2-C3	2.93	117.29	110.38
4	B	7066	T3A	C3-N1-C4	-2.90	111.92	116.17
5	A	600	FAD	C4X-C10-N10	2.87	120.59	116.48
5	B	600	FAD	O4'-C4'-C5'	-2.75	103.91	109.99
5	A	600	FAD	C4-C4X-C10	2.74	121.63	116.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	800	BOG	C1-O5-C5	-2.71	108.42	113.72
5	A	600	FAD	C2A-N1A-C6A	2.68	123.13	118.73
2	A	1949	BOG	C4-C3-C2	2.62	115.42	110.83
5	A	600	FAD	C4A-N9A-C8A	2.62	108.48	105.74
2	A	1949	BOG	C1-C2-C3	2.59	115.46	110.01
5	A	600	FAD	C8M-C8-C9	-2.49	115.18	119.57
6	A	700	PEP	O3P-P-O1P	-2.46	101.26	110.83
2	B	800	BOG	O2-C2-C1	2.46	115.93	110.08
5	B	600	FAD	C4-C4X-N5	2.43	121.56	118.21
5	B	600	FAD	N3A-C4A-N9A	2.41	131.26	127.17
6	B	700	PEP	O2P-P-O2	2.40	112.30	105.32
4	A	7066	T3A	O2-S1-O1	2.38	121.56	113.82
5	B	600	FAD	C10-C4X-N5	-2.37	119.97	124.81
4	B	7066	T3A	O2-S1-O1	2.36	121.50	113.82
6	A	700	PEP	O2-P-O1P	2.36	117.01	109.47
5	B	600	FAD	O2P-P-O3P	2.32	113.56	107.27
2	B	800	BOG	C1-C2-C3	-2.32	105.12	110.01
4	B	7066	T3A	O1-S1-C1	2.32	110.23	106.73
4	B	7066	T3A	O3-S1-O2	-2.31	105.63	111.40
6	B	700	PEP	O2P-P-O1P	-2.26	102.04	110.83
4	A	7066	T3A	O3-S1-O2	-2.26	105.75	111.40
5	A	600	FAD	O3B-C3B-C2B	2.25	119.05	111.82
6	B	700	PEP	O3P-P-O2P	2.25	116.25	107.80
4	A	7066	T3A	O1-S1-C1	2.18	110.02	106.73
5	A	600	FAD	C5X-N5-C4X	2.15	121.56	118.09
5	A	600	FAD	O3P-PA-O1A	-2.14	104.26	110.70
5	B	600	FAD	C9A-C5X-N5	-2.13	120.19	122.45
4	B	7066	T3A	O3-S1-C1	-2.13	101.84	106.00
5	B	600	FAD	C4X-C4-N3	2.12	118.66	113.25
5	B	600	FAD	O3B-C3B-C2B	2.11	118.59	111.82
2	A	800	BOG	O6-C6-C5	2.11	118.51	111.33
5	B	600	FAD	C4X-C10-N10	2.08	119.46	116.48
5	B	600	FAD	O4-C4-N3	-2.07	116.22	120.11
2	A	1949	BOG	O2-C2-C3	-2.07	105.50	110.38
4	A	7066	T3A	O3-S1-C1	-2.06	101.97	106.00
5	A	600	FAD	O3B-C3B-C4B	-2.03	105.26	111.08
5	A	600	FAD	C9A-C5X-N5	-2.01	120.32	122.45
2	B	800	BOG	C6-C5-C4	2.01	117.94	113.02
5	A	600	FAD	O4-C4-C4X	-2.00	121.25	126.53

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	800	BOG	C4
2	A	800	BOG	C3
2	A	1949	BOG	C3
2	A	1949	BOG	C2

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	700	PEP	O1-C1-C2-O2
6	A	700	PEP	O2'-C1-C2-O2
6	A	700	PEP	C3-C2-O2-P
6	A	700	PEP	C2-O2-P-O1P
6	B	700	PEP	O1-C1-C2-O2
6	B	700	PEP	O2'-C1-C2-O2
6	B	700	PEP	C3-C2-O2-P
2	A	900	BOG	O5-C1-O1-C1'
2	A	900	BOG	C2-C1-O1-C1'
2	A	900	BOG	O5-C5-C6-O6
2	B	800	BOG	O5-C1-O1-C1'
2	A	900	BOG	O1-C1'-C2'-C3'
2	B	800	BOG	O5-C5-C6-O6
2	A	800	BOG	O5-C5-C6-O6
2	B	900	BOG	O1-C1'-C2'-C3'
2	A	900	BOG	C1'-C2'-C3'-C4'
2	A	1949	BOG	C1'-C2'-C3'-C4'
2	A	800	BOG	C3'-C4'-C5'-C6'
2	A	1949	BOG	C4'-C5'-C6'-C7'
3	A	1954	EDO	O1-C1-C2-O2
3	A	7069	EDO	O1-C1-C2-O2
2	A	1949	BOG	C3'-C4'-C5'-C6'
2	A	1949	BOG	C2'-C3'-C4'-C5'
3	A	1953	EDO	O1-C1-C2-O2
3	A	1955	EDO	O1-C1-C2-O2
3	B	901	EDO	O1-C1-C2-O2
3	B	902	EDO	O1-C1-C2-O2
3	B	903	EDO	O1-C1-C2-O2
3	B	907	EDO	O1-C1-C2-O2
3	B	7069	EDO	O1-C1-C2-O2
2	B	800	BOG	O1-C1'-C2'-C3'
2	A	900	BOG	C4-C5-C6-O6
2	A	800	BOG	C5'-C6'-C7'-C8'
2	B	800	BOG	C4-C5-C6-O6
3	A	1950	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	1956	EDO	O1-C1-C2-O2
2	B	800	BOG	C1'-C2'-C3'-C4'
2	B	900	BOG	C4-C5-C6-O6
5	A	600	FAD	N10-C1'-C2'-O2'
5	B	600	FAD	N10-C1'-C2'-O2'
3	B	904	EDO	O1-C1-C2-O2
2	A	1949	BOG	C2'-C1'-O1-C1
3	A	1951	EDO	O1-C1-C2-O2
3	A	1952	EDO	O1-C1-C2-O2
3	B	906	EDO	O1-C1-C2-O2
6	B	700	PEP	C2-O2-P-O3P
3	A	7070	EDO	O1-C1-C2-O2
3	B	7067	EDO	O1-C1-C2-O2
2	B	800	BOG	C5'-C6'-C7'-C8'
2	B	900	BOG	C2'-C3'-C4'-C5'
5	A	600	FAD	O4B-C4B-C5B-O5B

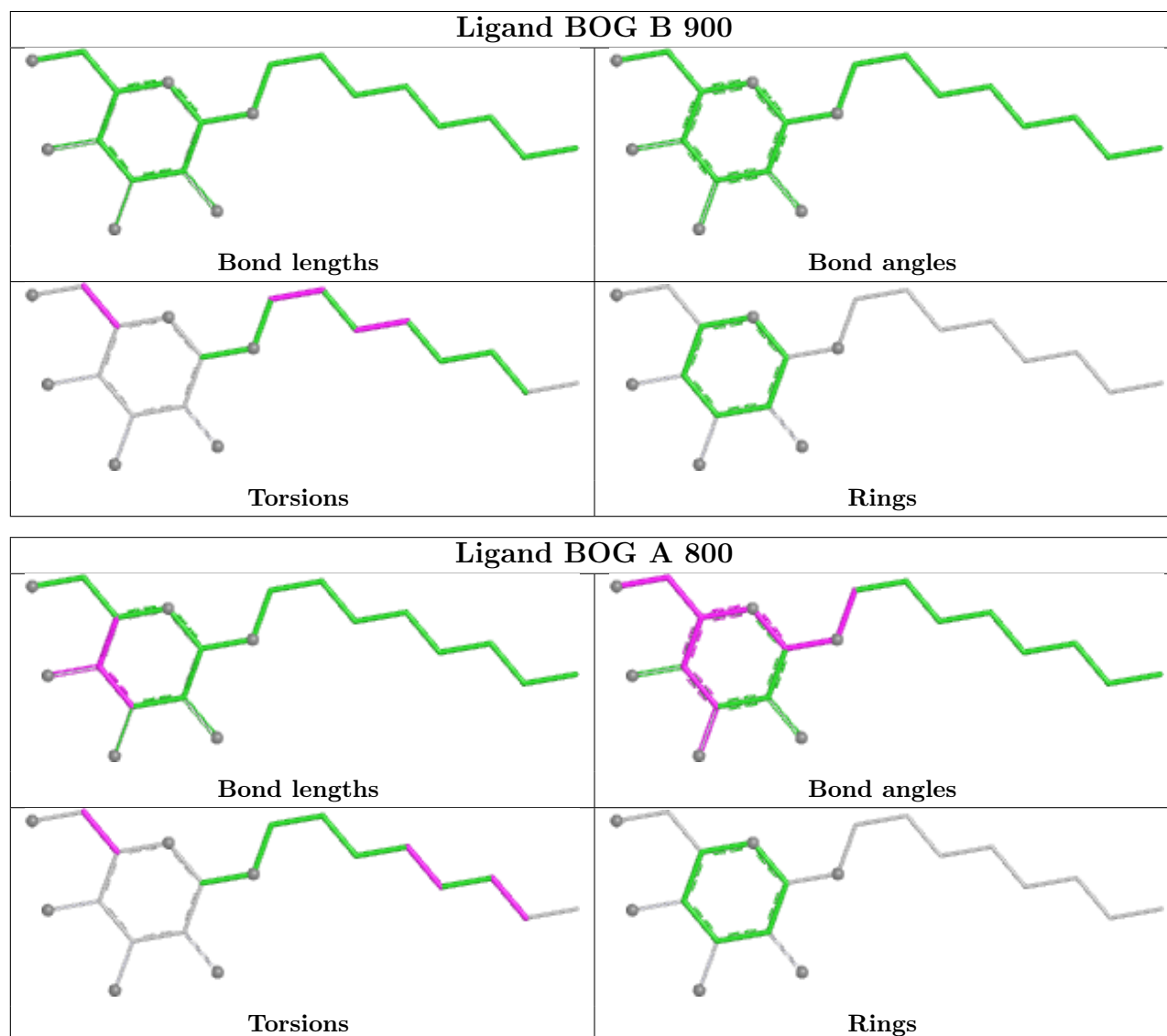
There are no ring outliers.

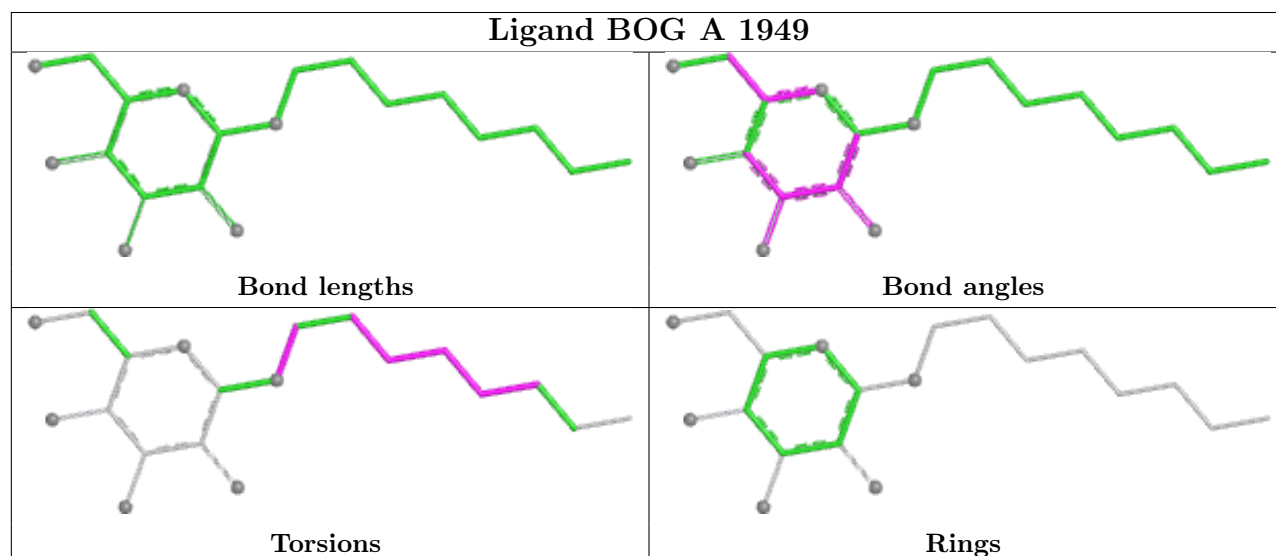
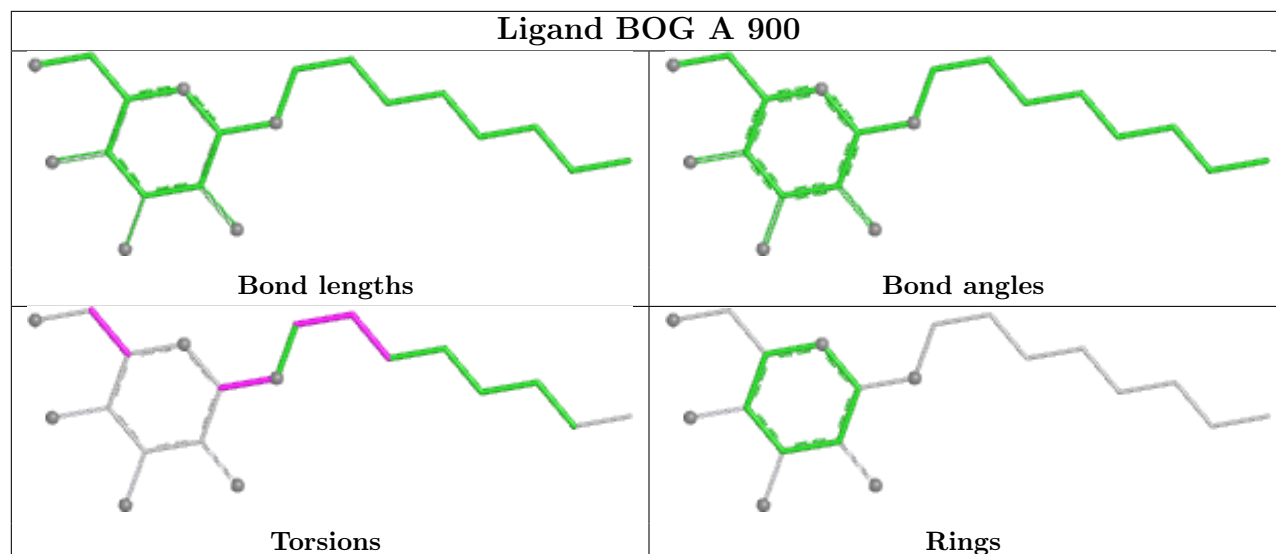
17 monomers are involved in 53 short contacts:

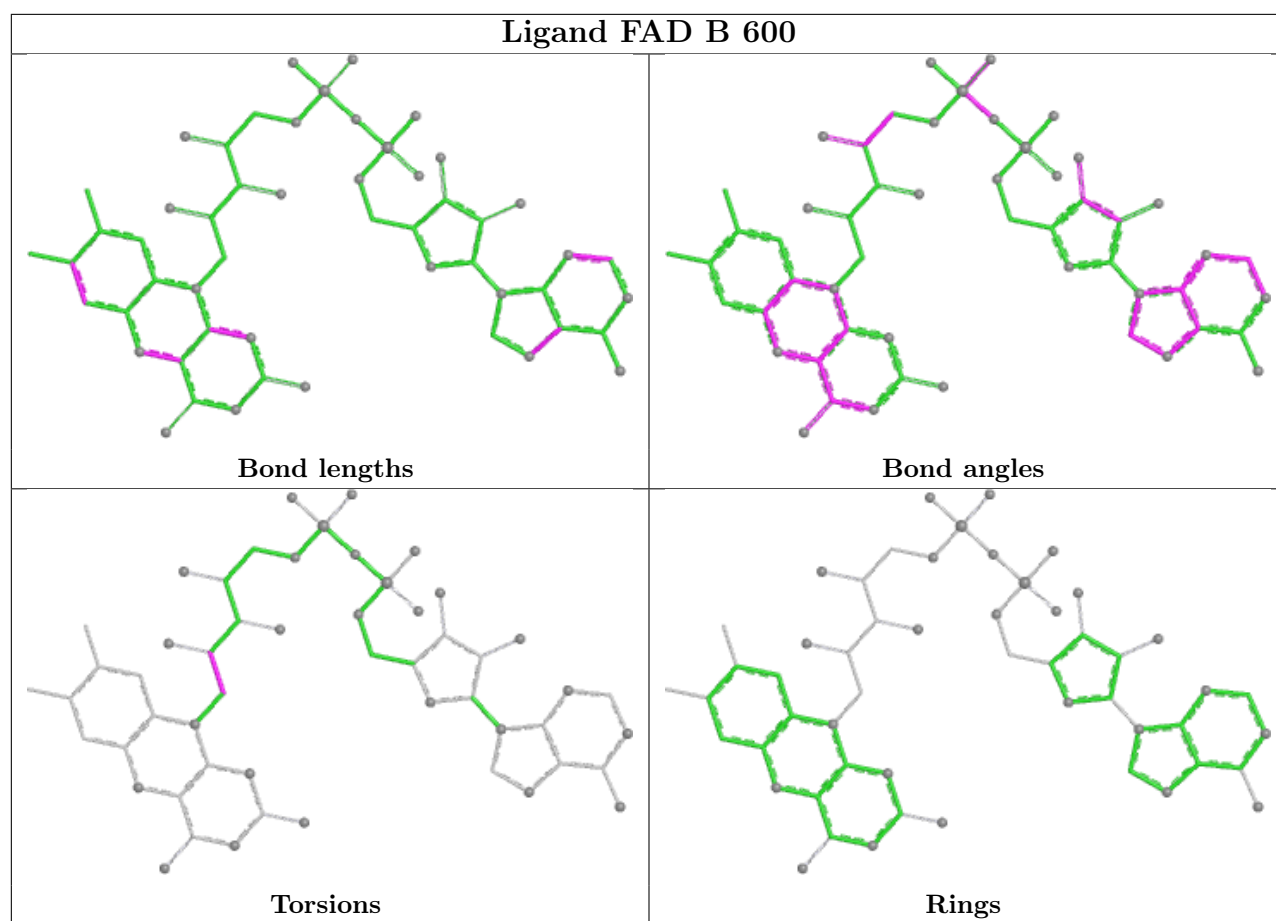
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1953	EDO	1	0
7	B	7068	IMD	2	0
3	A	7069	EDO	1	0
6	A	700	PEP	4	0
2	B	900	BOG	7	0
2	A	800	BOG	16	0
2	A	900	BOG	2	0
3	A	1956	EDO	1	0
3	B	7067	EDO	2	0
2	A	1949	BOG	4	0
6	B	700	PEP	3	0
5	B	600	FAD	1	0
4	B	7066	T3A	1	0
5	A	600	FAD	1	0
2	B	800	BOG	3	0
3	B	901	EDO	2	0
3	A	1954	EDO	2	0

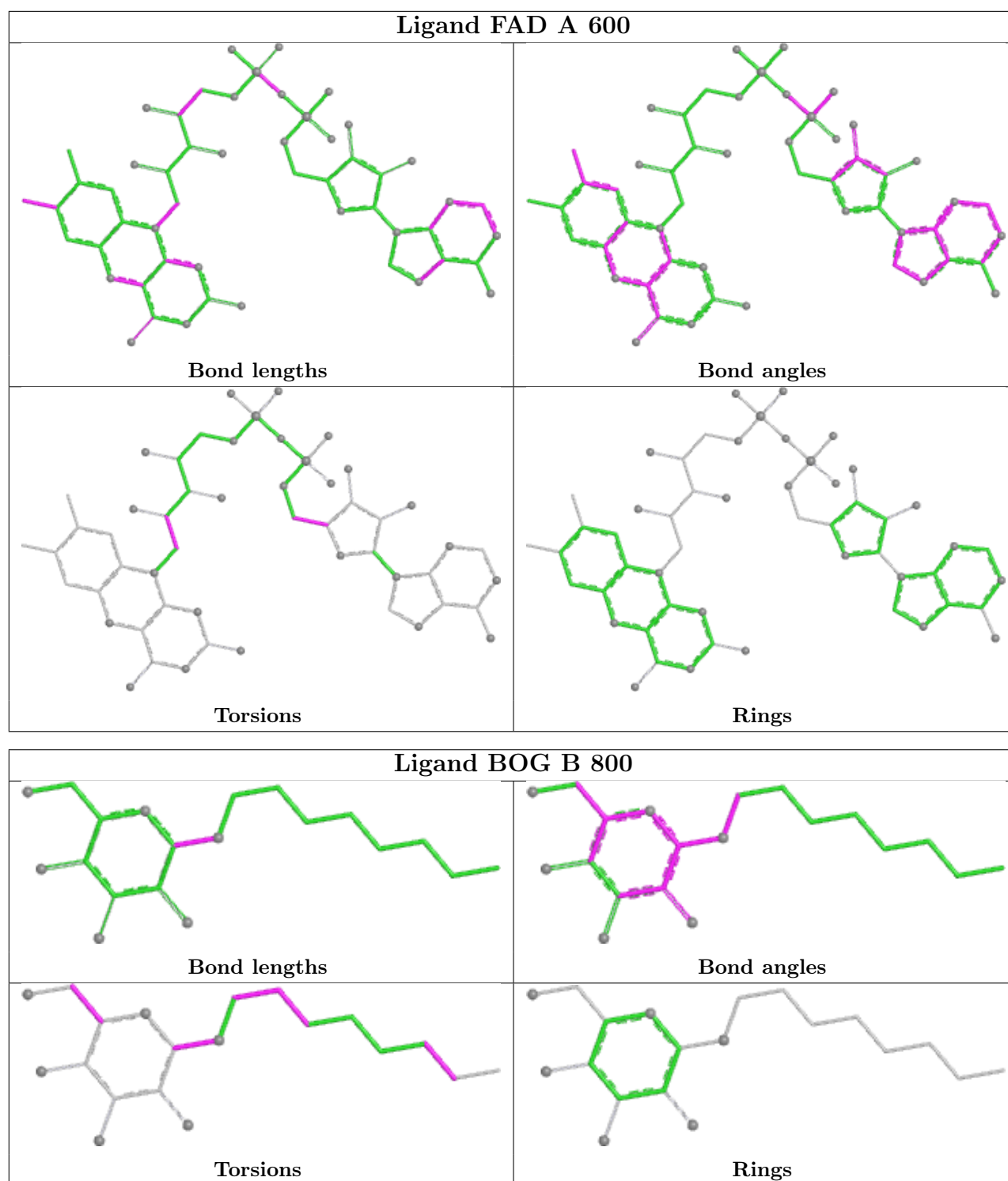
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

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 ⓘ

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	495/501 (98%)	-1.49	0 100 100	15, 32, 52, 81	0
1	B	495/501 (98%)	-1.48	0 100 100	16, 32, 52, 80	0
All	All	990/1002 (98%)	-1.49	0 100 100	15, 32, 52, 81	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BOG	A	900	20/20	0.97	0.12	95,98,102,102	0
2	BOG	A	800	20/20	0.97	0.08	106,111,112,112	0
2	BOG	B	900	20/20	0.97	0.12	95,97,100,101	0
2	BOG	A	1949	20/20	0.98	0.08	80,86,90,91	0
2	BOG	B	800	20/20	0.98	0.07	100,101,104,105	0
3	EDO	A	1950	4/4	0.98	0.07	51,54,56,58	0
3	EDO	A	1951	4/4	0.98	0.05	26,38,39,42	0

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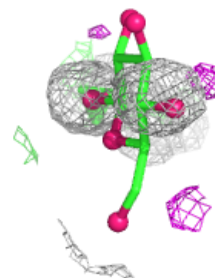
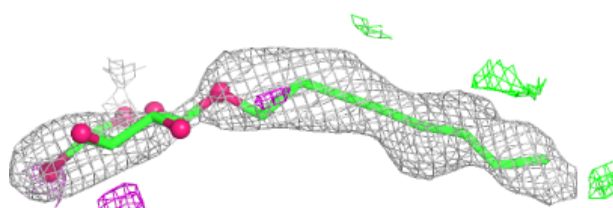
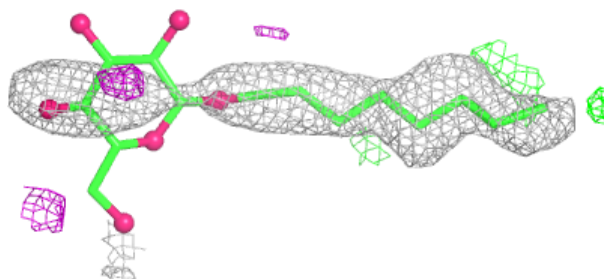
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	A	1953	4/4	0.98	0.06	50,50,51,52	0
3	EDO	A	1954	4/4	0.98	0.05	35,39,40,40	0
3	EDO	A	1956	4/4	0.98	0.06	50,51,52,54	0
3	EDO	A	7069	4/4	0.98	0.05	52,53,53,54	0
3	EDO	B	901	4/4	0.98	0.06	38,41,44,46	0
3	EDO	B	902	4/4	0.98	0.08	54,56,57,59	0
3	EDO	B	904	4/4	0.98	0.05	25,37,41,41	0
3	EDO	A	7068	4/4	0.99	0.07	62,64,65,65	0
3	EDO	A	1955	4/4	0.99	0.07	64,66,68,69	0
3	EDO	A	1952	4/4	0.99	0.03	47,47,48,51	0
3	EDO	A	1957	4/4	0.99	0.05	57,60,61,63	0
3	EDO	B	903	4/4	0.99	0.03	51,51,52,55	0
3	EDO	A	7067	4/4	0.99	0.06	60,63,64,65	0
3	EDO	B	905	4/4	0.99	0.05	56,58,58,58	0
3	EDO	B	906	4/4	0.99	0.07	54,57,58,58	0
3	EDO	B	907	4/4	0.99	0.04	60,61,62,62	0
3	EDO	B	908	4/4	0.99	0.08	44,44,46,47	0
3	EDO	B	7067	4/4	0.99	0.03	48,52,52,55	0
3	EDO	B	7069	4/4	0.99	0.03	20,20,20,20	0
4	T3A	A	7066	15/15	0.99	0.03	26,36,44,45	0
4	T3A	B	7066	15/15	0.99	0.03	25,37,44,46	0
6	PEP	A	700	10/10	0.99	0.03	30,34,39,43	0
7	IMD	B	7068	5/5	0.99	0.04	32,44,45,47	0
5	FAD	B	600	53/53	1.00	0.02	15,19,21,22	0
3	EDO	A	7070	4/4	1.00	0.02	20,20,20,20	0
6	PEP	B	700	10/10	1.00	0.03	25,31,38,39	0
5	FAD	A	600	53/53	1.00	0.02	14,18,21,22	0

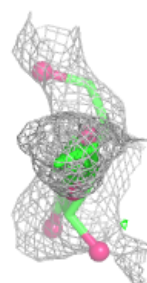
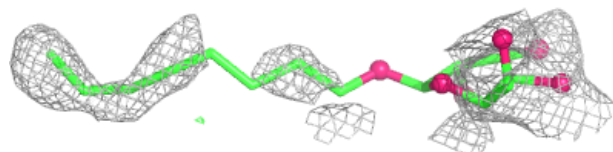
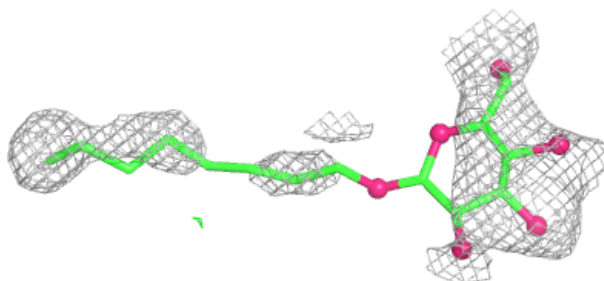
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BOG A 900:

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

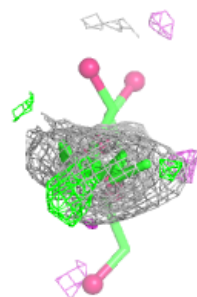
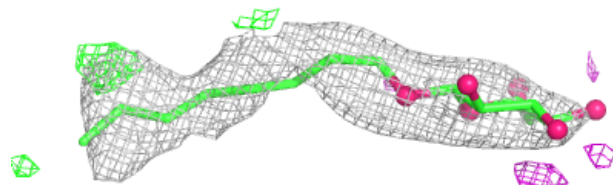
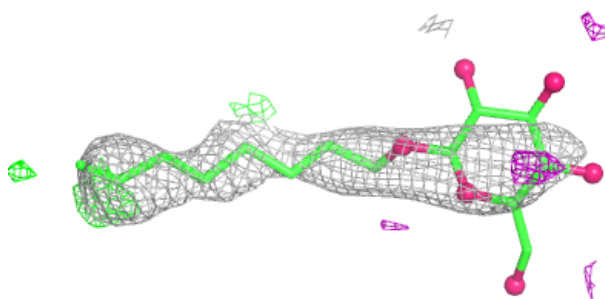
**Electron density around BOG A 800:**

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



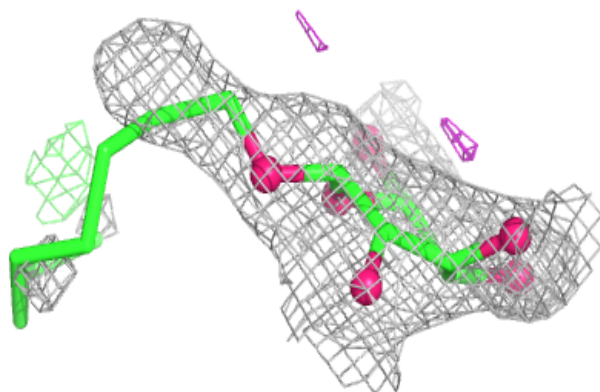
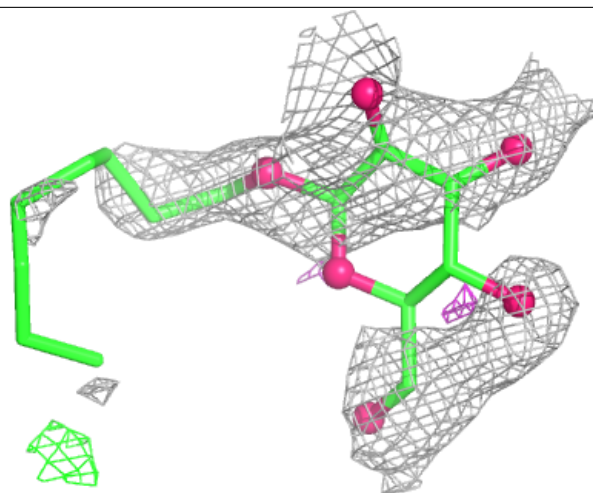
Electron density around BOG B 900:

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



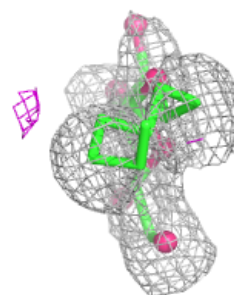
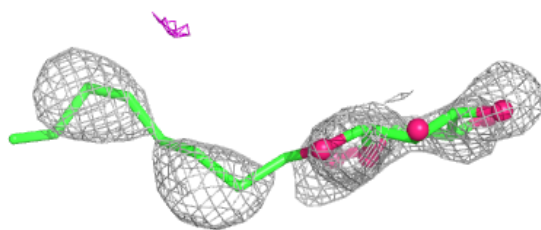
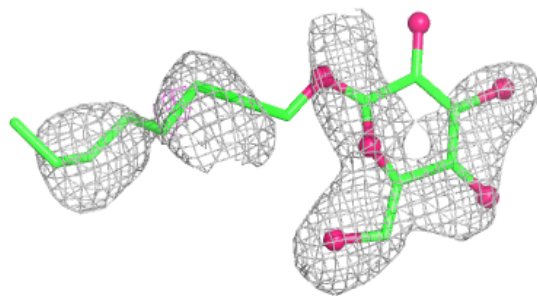
Electron density around BOG A 1949:

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

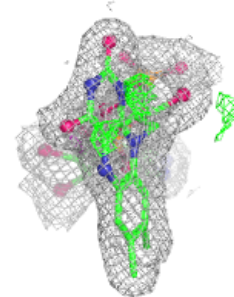
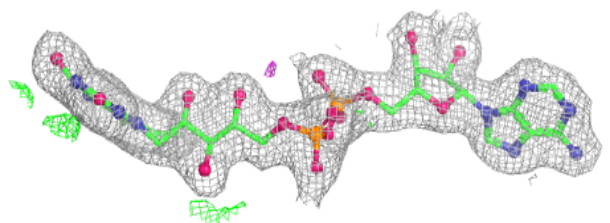
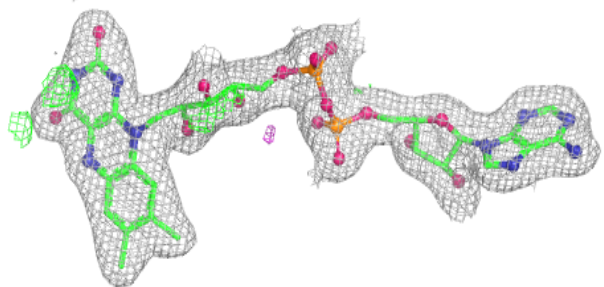


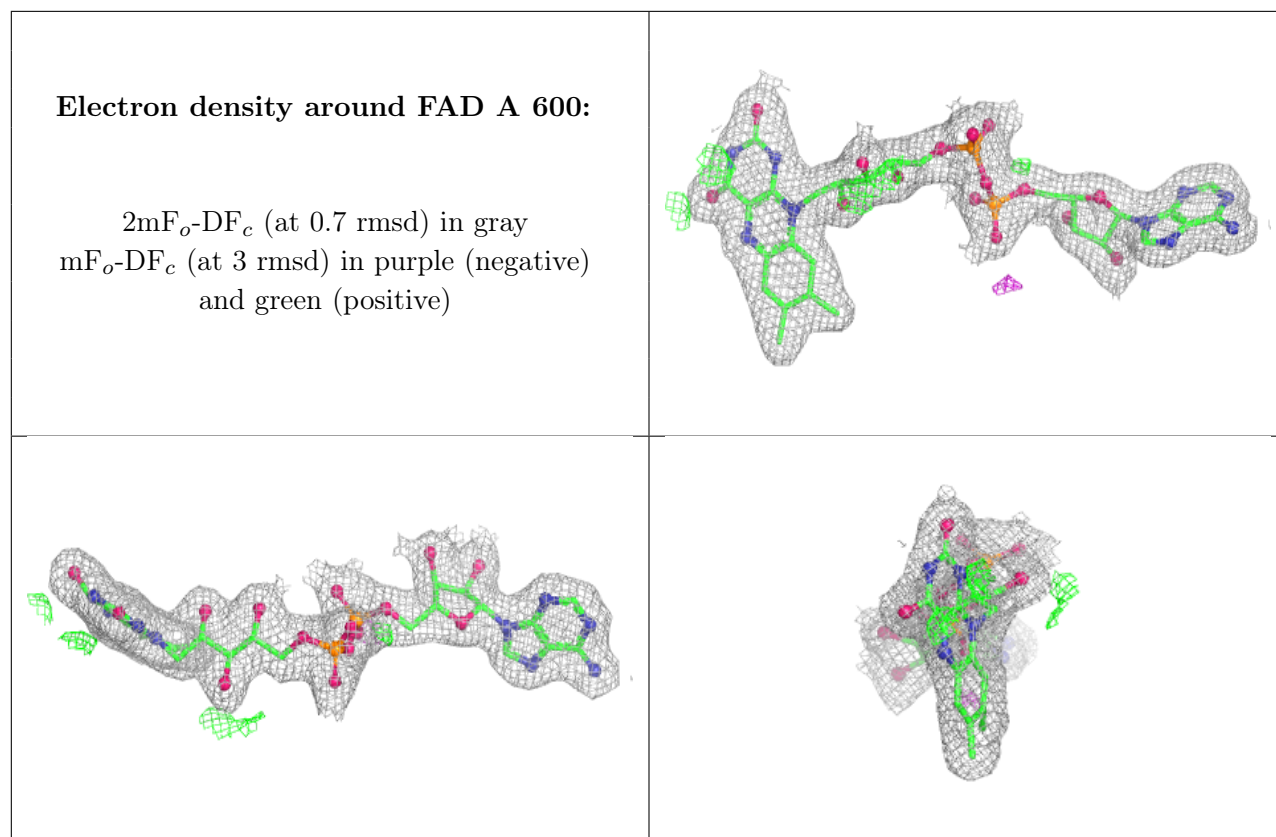
Electron density around BOG B 800:

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

**Electron density around FAD B 600:**

$2mF_o-DF_c$ (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.