Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:49 AM UTC

PDB ID	:	7BZV / pdb_00007bzv
Title	:	Crystal structure of 2-aminomuconic 6-semialdehyde dehydrogenase from Pseudomonas species AP-3
Authors	:	Shi, Q.L.; Chen, Y.J.; Su, D.
Deposited on	:	2020-04-28
Resolution	:	1.99 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

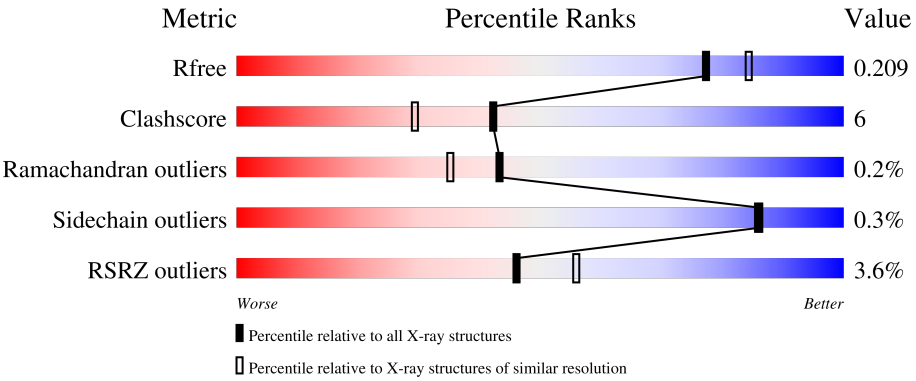
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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

Mol	Chain	Length	Quality of chain
1	A	492	3% 89% 9%
1	B	492	4% 89% 10%

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	PEG	A	504	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15883 atoms, of which 7528 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.

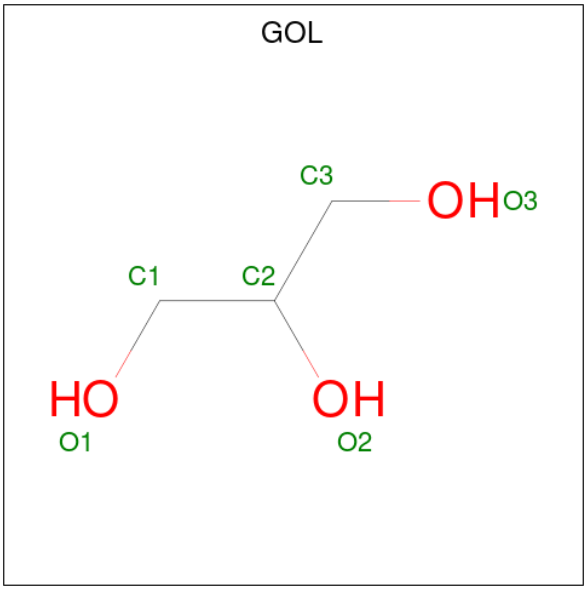
- Molecule 1 is a protein called 2-aminomuconic 6-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	489	Total	C	H	N	O	S	0	2	0
			7501	2380	3726	671	705	19			
1	B	490	Total	C	H	N	O	S	0	2	0
			7515	2385	3733	672	706	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q9KWS5
B	0	SER	-	expression tag	UNP Q9KWS5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



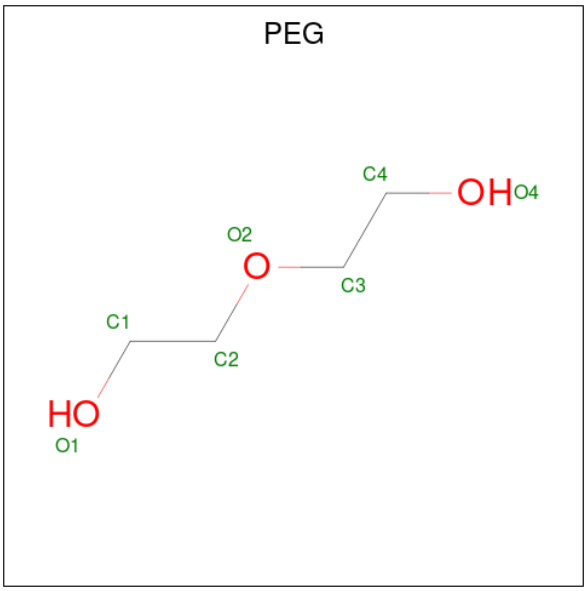
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			9	3	3	3		

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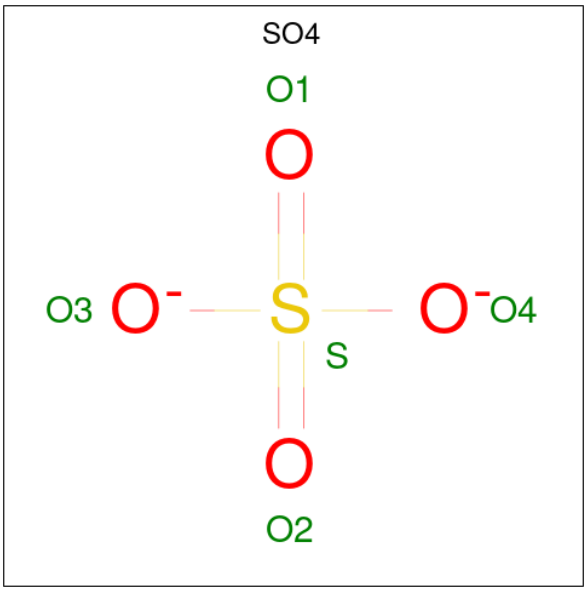
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C H O 14 3 8 3	0	0
2	B	1	Total C H O 14 3 8 3	0	0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 17 4 10 3	0	0
3	A	1	Total C H O 17 4 10 3	0	0
3	A	1	Total C H O 17 4 10 3	0	0
3	B	1	Total C H O 17 4 10 3	0	0
3	B	1	Total C H O 17 4 10 3	0	0

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is water.

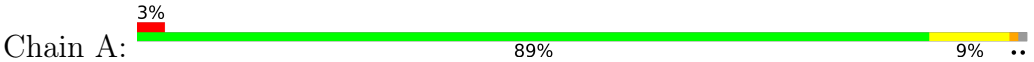
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	402	Total O 402 402	0	0
5	B	338	Total O 338 338	0	0

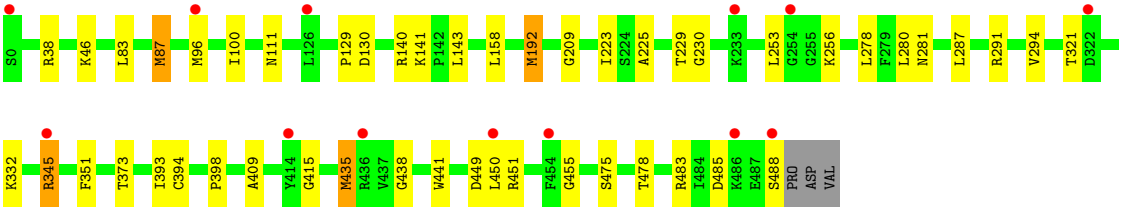
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 2-aminomuconic 6-semialdehyde dehydrogenase

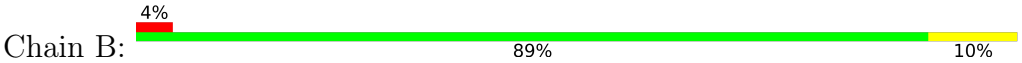
Chain A:

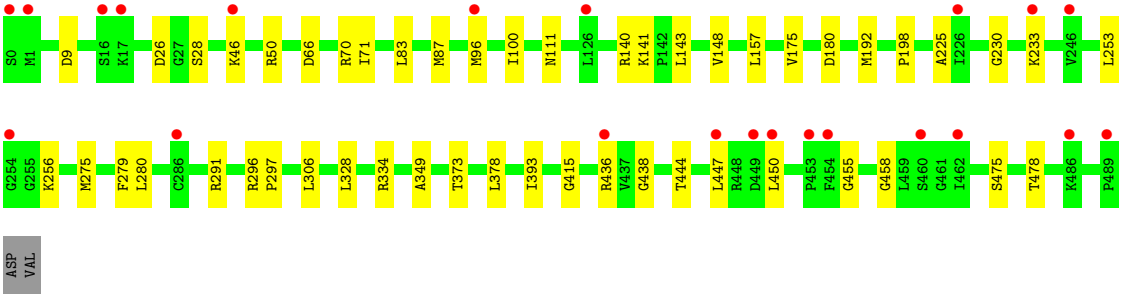




• Molecule 1: 2-aminomuconic 6-semialdehyde dehydrogenase

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	181.53Å 181.53Å 181.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.54 – 1.99 25.54 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.8 (25.54-1.99) 99.7 (25.54-1.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.09 (at 1.98Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.172 , 0.207 0.173 , 0.209	Depositor DCC
R_{free} test set	6091 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 57.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	15883	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GOL, PEG

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	0.58	2/3867 (0.1%)	0.75	3/5249 (0.1%)
1	B	0.51	1/3877 (0.0%)	0.68	0/5263
All	All	0.54	3/7744 (0.0%)	0.72	3/10512 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	192	MET	SD-CE	9.44	2.03	1.79
1	A	87	MET	SD-CE	-5.97	1.64	1.79
1	B	198	PRO	CA-C	5.00	1.54	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	435	MET	CG-SD-CE	11.73	126.71	100.90
1	A	321	THR	CA-C-N	5.82	130.57	120.68
1	A	321	THR	C-N-CA	5.82	130.57	120.68

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	3775	3726	3717	45	0
1	B	3782	3733	3719	48	0
2	A	6	3	8	1	0
2	B	12	16	16	1	0
3	A	21	30	30	7	0
3	B	14	20	20	1	0
4	B	5	0	0	0	0
5	A	402	0	0	6	0
5	B	338	0	0	6	1
All	All	8355	7528	7510	95	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (95) 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:192:MET:SD	1:A:192:MET:CE	2.03	1.47
1:A:409:ALA:HB1	1:A:435:MET:HE1	1.18	1.16
1:B:275:MET:HE3	1:B:306:LEU:CD1	1.78	1.11
1:A:83:LEU:HG	1:A:87:MET:HE2	1.32	1.11
1:B:275:MET:HE3	1:B:306:LEU:HD11	1.17	1.10
1:A:409:ALA:CB	1:A:435:MET:HE1	1.82	1.08
1:A:435:MET:HE3	5:A:885:HOH:O	1.51	1.07
1:A:409:ALA:HB1	1:A:435:MET:CE	1.95	0.96
1:B:141:LYS:NZ	1:B:478:THR:OG1	2.02	0.92
1:B:275:MET:CE	1:B:306:LEU:HD11	2.01	0.89
1:A:96:MET:HE2	1:A:281:ASN:HB2	1.54	0.88
1:A:111:ASN:CG	1:A:450:LEU:HD12	1.99	0.86
1:B:83:LEU:HD22	1:B:87:MET:HE2	1.56	0.86
1:B:233:LYS:H	1:B:233:LYS:HE2	1.43	0.82
1:B:96:MET:SD	1:B:280:LEU:HD11	2.19	0.81
1:A:111:ASN:OD1	3:A:504:PEG:H31	1.82	0.79
1:A:83:LEU:HG	1:A:87:MET:CE	2.12	0.78
1:A:111:ASN:OD1	1:A:450:LEU:HD12	1.83	0.76
1:B:296:ARG:NH2	5:B:601:HOH:O	2.14	0.76
1:B:26:ASP:OD1	1:B:28:SER:OG	2.05	0.75
1:A:141:LYS:HE2	5:A:894:HOH:O	1.88	0.73
1:A:449:ASP:HA	3:A:504:PEG:H42	1.70	0.72
1:A:96:MET:HE1	1:A:280:LEU:HG	1.77	0.66
1:A:87:MET:HE3	5:A:618:HOH:O	1.94	0.66
1:B:96:MET:HE1	1:B:280:LEU:CD2	2.24	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:LYS:HE2	5:B:874:HOH:O	1.95	0.65
1:A:38:ARG:NH1	5:A:601:HOH:O	2.30	0.64
1:B:96:MET:SD	1:B:280:LEU:CD1	2.86	0.63
1:B:96:MET:HE1	1:B:280:LEU:HD21	1.81	0.62
1:A:87:MET:CE	5:A:618:HOH:O	2.48	0.61
1:A:96:MET:SD	1:A:280:LEU:HD11	2.40	0.61
1:B:111:ASN:OD1	1:B:450:LEU:HD12	2.01	0.61
1:B:275:MET:HE3	1:B:306:LEU:HD12	1.77	0.61
1:B:141:LYS:NZ	1:B:478:THR:CB	2.66	0.58
1:B:87:MET:HE3	5:B:609:HOH:O	2.03	0.57
1:A:483:ARG:NH1	1:A:485:ASP:OD1	2.37	0.57
3:A:504:PEG:O1	3:A:504:PEG:H32	2.05	0.56
1:B:148:VAL:HB	1:B:175:VAL:HG22	1.88	0.56
1:A:192:MET:CE	1:A:192:MET:CG	2.83	0.56
1:A:96:MET:CE	1:A:280:LEU:HG	2.36	0.55
1:A:345[B]:ARG:HG2	1:A:351:PHE:CE1	2.41	0.55
1:B:275:MET:HE1	1:B:279:PHE:HE2	1.71	0.55
1:A:451:ARG:HD3	3:A:504:PEG:H21	1.88	0.54
1:A:451:ARG:CD	3:A:504:PEG:H21	2.38	0.53
1:B:180:ASP:OD2	2:B:504:GOL:H32	2.10	0.52
1:B:233:LYS:H	1:B:233:LYS:CE	2.20	0.52
1:A:140:ARG:HG3	1:A:475:SER:OG	2.12	0.50
1:B:140:ARG:HG3	1:B:475:SER:OG	2.13	0.49
1:A:96:MET:SD	1:A:280:LEU:CD1	3.00	0.49
1:A:158:LEU:HD13	1:A:229:THR:HG21	1.94	0.49
1:B:143:LEU:HD13	1:B:225:ALA:HB2	1.94	0.49
1:A:96:MET:HE1	1:A:280:LEU:CD2	2.43	0.49
3:B:505:PEG:H22	5:B:845:HOH:O	2.13	0.49
1:B:275:MET:CE	1:B:306:LEU:CD1	2.70	0.48
1:B:66:ASP:HB3	5:B:865:HOH:O	2.13	0.48
1:B:157:LEU:C	1:B:157:LEU:HD23	2.39	0.48
1:B:296:ARG:HB3	1:B:297:PRO:HD3	1.96	0.47
1:B:46:LYS:HE3	1:B:46:LYS:HA	1.96	0.47
5:A:894:HOH:O	1:B:436:ARG:HG3	2.15	0.47
1:A:46:LYS:HB3	1:A:46:LYS:NZ	2.30	0.46
1:B:458:GLY:O	5:B:602:HOH:O	2.20	0.46
1:A:96:MET:HE1	1:A:280:LEU:CG	2.45	0.46
1:B:9:ASP:OD2	1:B:50:ARG:NH1	2.49	0.46
1:B:140:ARG:HG3	1:B:475:SER:CB	2.46	0.46
1:A:450:LEU:H	3:A:504:PEG:H42	1.81	0.46
1:A:209:GLY:HA3	3:A:503:PEG:H42	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:MET:HE1	1:B:280:LEU:HG	1.98	0.45
1:B:349:ALA:HB2	1:B:378:LEU:HD11	1.98	0.45
1:A:278:LEU:HD22	1:A:394:CYS:HB3	1.98	0.45
1:B:373:THR:O	1:B:393:ILE:HA	2.16	0.45
1:A:141:LYS:NZ	1:A:478:THR:OG1	2.48	0.45
1:B:256:LYS:CG	1:B:291:ARG:HD2	2.47	0.45
1:B:275:MET:CE	1:B:279:PHE:HE2	2.29	0.45
1:A:100:ILE:HD12	1:A:280:LEU:HD21	1.99	0.44
1:B:438:GLY:HA3	1:B:455:GLY:O	2.18	0.44
1:A:129:PRO:HB2	1:A:130:ASP:OD2	2.17	0.44
1:A:373:THR:O	1:A:393:ILE:HA	2.18	0.43
1:A:256:LYS:HG3	1:A:291:ARG:CD	2.49	0.43
1:B:230:GLY:O	1:B:253:LEU:HA	2.19	0.43
1:B:96:MET:HE1	1:B:280:LEU:CG	2.48	0.43
1:A:223:ILE:HD12	1:A:223:ILE:HA	1.92	0.42
1:A:230:GLY:O	1:A:253:LEU:HA	2.20	0.42
1:B:328:LEU:HD12	1:B:334:ARG:HA	2.00	0.42
1:A:438:GLY:HA3	1:A:455:GLY:O	2.20	0.42
1:B:140:ARG:CG	1:B:475:SER:OG	2.67	0.42
1:B:100:ILE:CD1	1:B:280:LEU:HD21	2.50	0.42
1:A:332:LYS:HZ2	2:A:501:GOL:H12	1.85	0.42
1:B:71:ILE:HD13	1:B:192:MET:HE1	2.01	0.42
1:A:143:LEU:HD13	1:A:225:ALA:HB2	2.02	0.42
1:B:26:ASP:CG	1:B:28:SER:HG	2.13	0.41
1:A:287:LEU:HD13	1:A:441:TRP:HH2	1.86	0.41
1:B:96:MET:CE	1:B:280:LEU:HG	2.50	0.41
1:A:294:VAL:O	1:A:398:PRO:HA	2.22	0.40
1:B:66:ASP:O	1:B:70:ARG:HG3	2.21	0.40
1:B:444:THR:HG21	1:B:447:LEU:HD12	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:901:HOH:O	5:B:901:HOH:O[10_665]	2.10	0.10

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	489/492 (99%)	481 (98%)	7 (1%)	1 (0%)	43 35
1	B	490/492 (100%)	475 (97%)	14 (3%)	1 (0%)	43 35
All	All	979/984 (100%)	956 (98%)	21 (2%)	2 (0%)	43 35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	415	GLY
1	A	415	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	398/399 (100%)	395 (99%)	3 (1%)	73 71
1	B	399/399 (100%)	399 (100%)	0	100 100
All	All	797/798 (100%)	794 (100%)	3 (0%)	86 82

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	345[A]	ARG
1	A	345[B]	ARG
1	A	488	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

9 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	PEG	A	503	-	6,6,6	0.53	0	5,5,5	0.50	0
3	PEG	A	502	-	6,6,6	0.54	0	5,5,5	0.42	0
2	GOL	B	503	-	5,5,5	0.80	0	5,5,5	0.82	0
3	PEG	A	504	-	6,6,6	0.46	0	5,5,5	0.37	0
4	SO4	B	502	-	4,4,4	0.21	0	6,6,6	0.15	0
2	GOL	B	504	-	5,5,5	1.21	0	5,5,5	1.51	1 (20%)
3	PEG	B	501	-	6,6,6	0.51	0	5,5,5	0.34	0
2	GOL	A	501	-	5,5,5	1.14	1 (20%)	5,5,5	0.83	0
3	PEG	B	505	-	6,6,6	0.43	0	5,5,5	0.25	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	PEG	A	503	-	-	2/4/4/4	-
3	PEG	A	502	-	-	1/4/4/4	-
2	GOL	B	503	-	-	2/4/4/4	-
3	PEG	A	504	-	-	3/4/4/4	-
2	GOL	B	504	-	-	0/4/4/4	-
3	PEG	B	501	-	-	3/4/4/4	-
2	GOL	A	501	-	-	4/4/4/4	-
3	PEG	B	505	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	GOL	O2-C2	-2.07	1.37	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	504	GOL	C3-C2-C1	-2.64	102.13	111.80

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	501	PEG	O2-C3-C4-O4
3	A	503	PEG	O1-C1-C2-O2
3	B	501	PEG	O1-C1-C2-O2
2	A	501	GOL	O1-C1-C2-C3
2	A	501	GOL	C1-C2-C3-O3
2	B	503	GOL	C1-C2-C3-O3
3	A	502	PEG	O2-C3-C4-O4
2	A	501	GOL	O1-C1-C2-O2
2	B	503	GOL	O2-C2-C3-O3
3	A	503	PEG	O2-C3-C4-O4
3	A	504	PEG	O2-C3-C4-O4
3	A	504	PEG	C4-C3-O2-C2
3	B	501	PEG	C1-C2-O2-C3
3	B	505	PEG	O1-C1-C2-O2
3	B	505	PEG	C4-C3-O2-C2
3	A	504	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
2	A	501	GOL	O2-C2-C3-O3

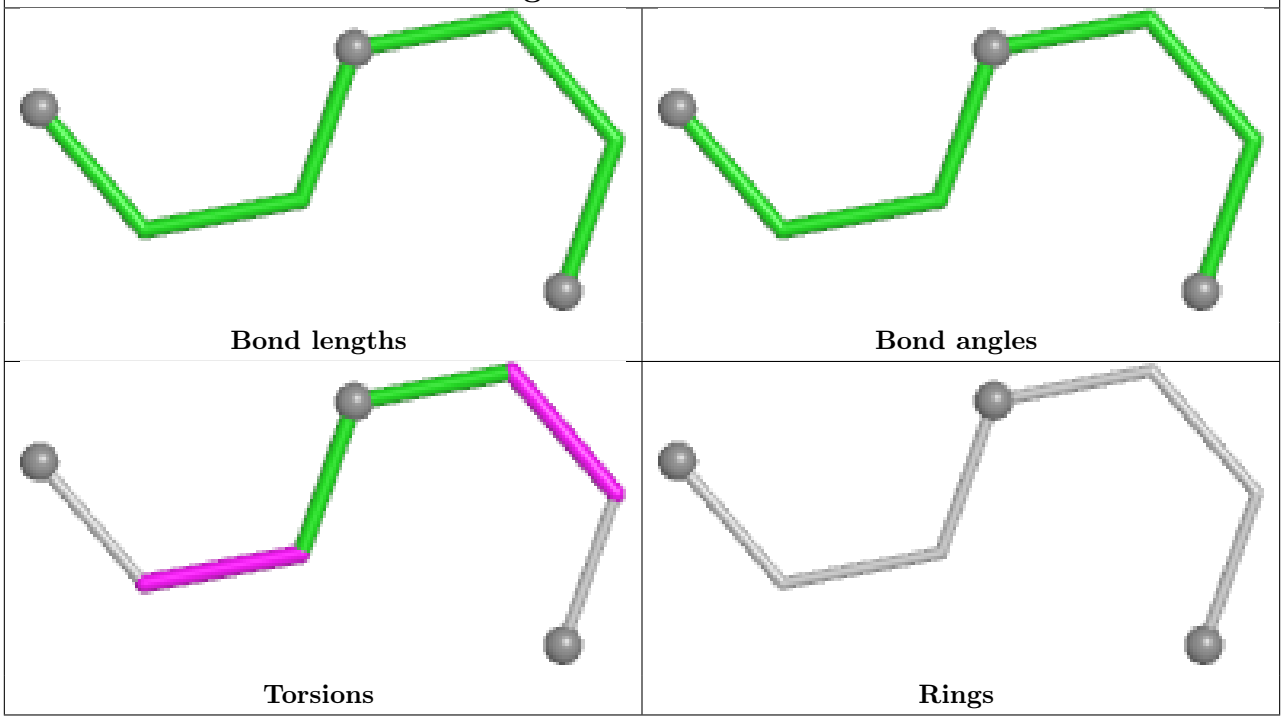
There are no ring outliers.

5 monomers are involved in 10 short contacts:

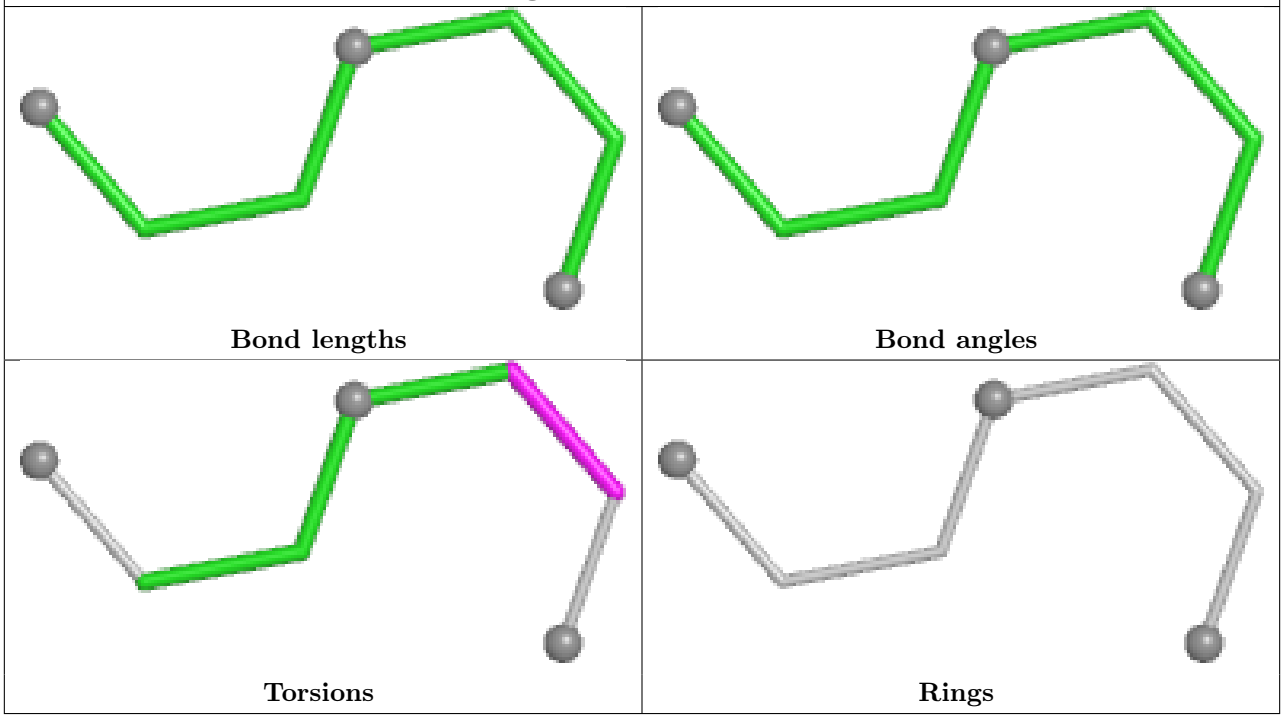
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	PEG	1	0
3	A	504	PEG	6	0
2	B	504	GOL	1	0
2	A	501	GOL	1	0
3	B	505	PEG	1	0

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

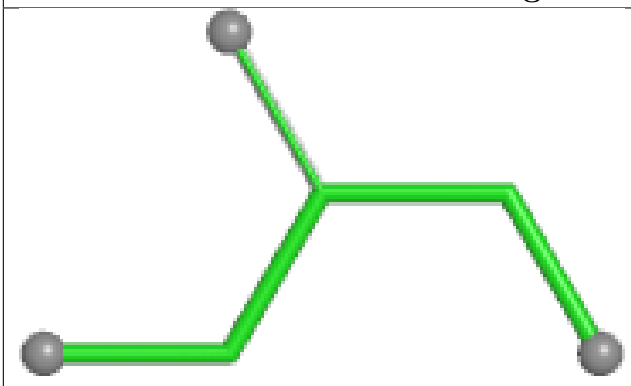
Ligand PEG A 503



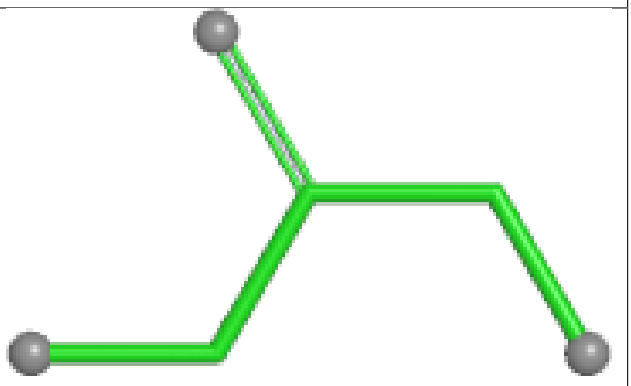
Ligand PEG A 502



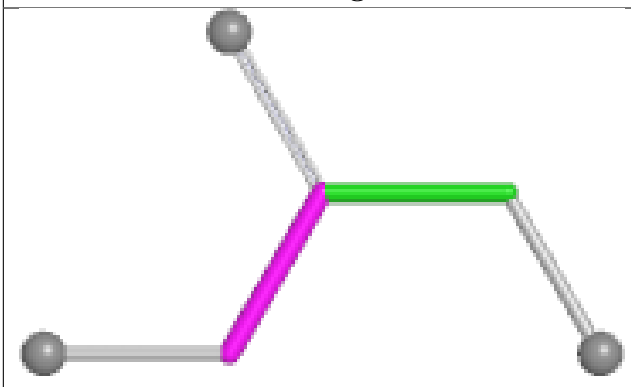
Ligand GOL B 503



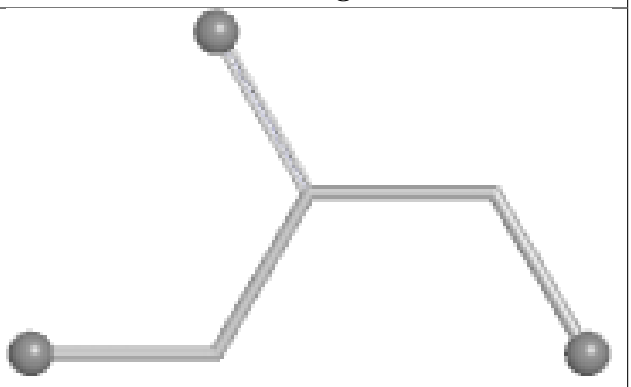
Bond lengths



Bond angles

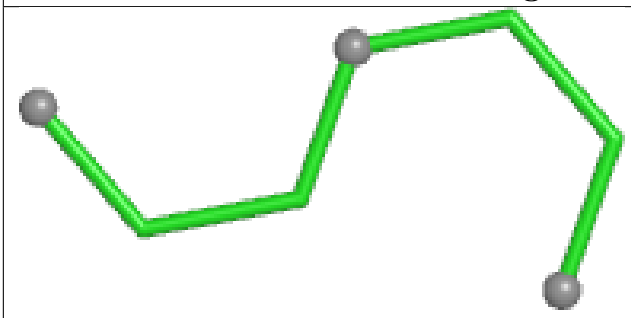


Torsions

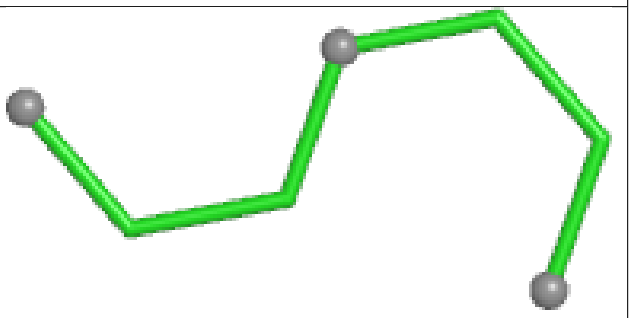


Rings

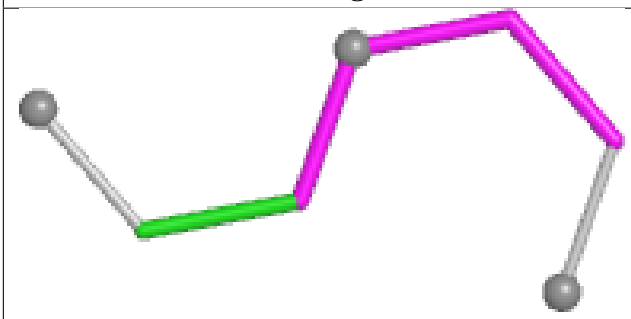
Ligand PEG A 504



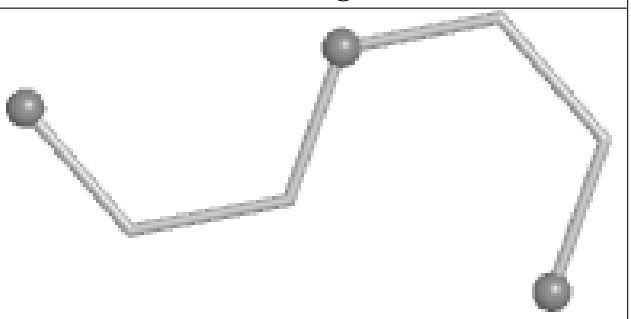
Bond lengths



Bond angles

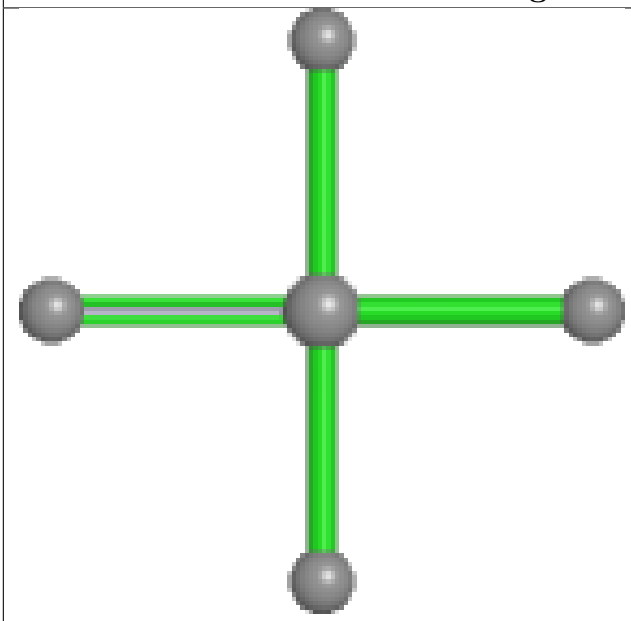


Torsions

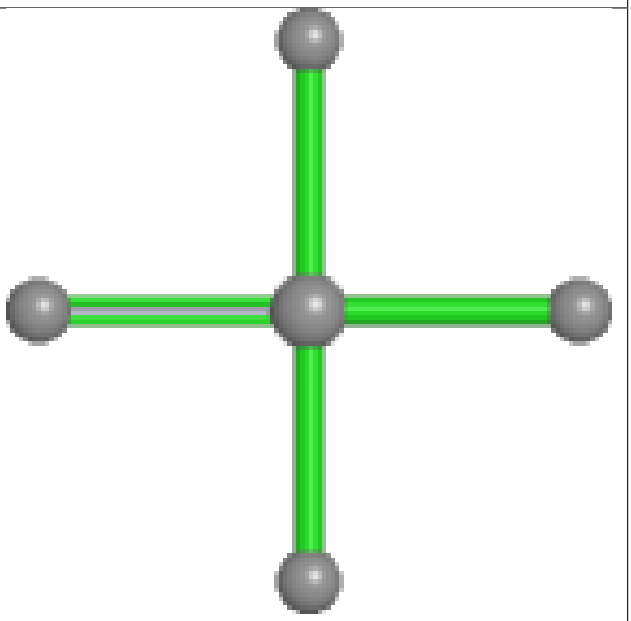


Rings

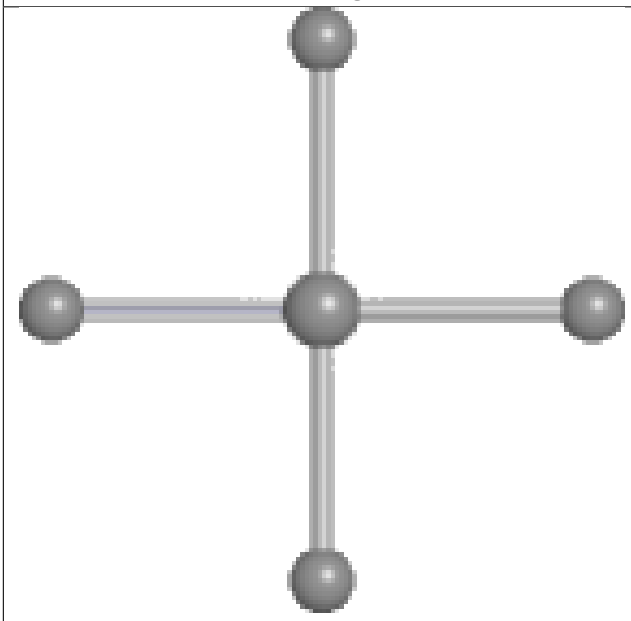
Ligand SO4 B 502



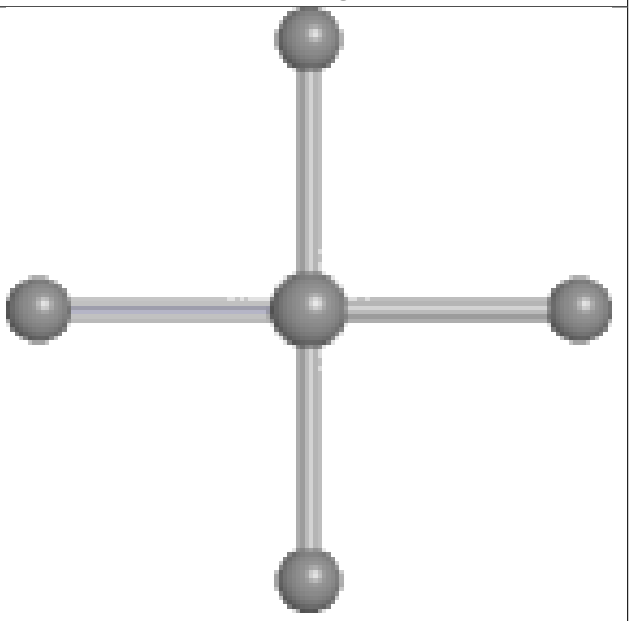
Bond lengths



Bond angles

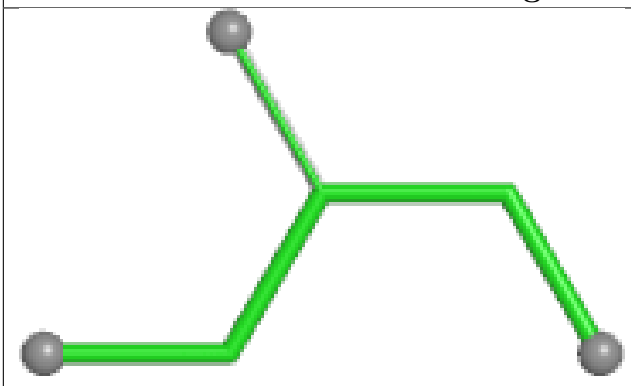


Torsions

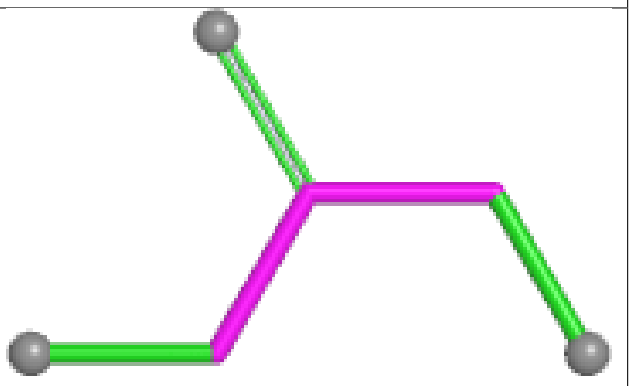


Rings

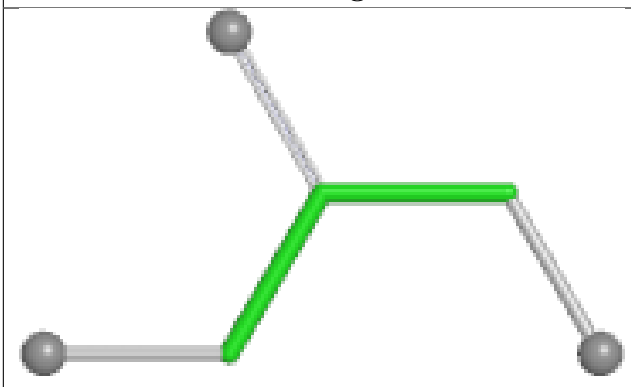
Ligand GOL B 504



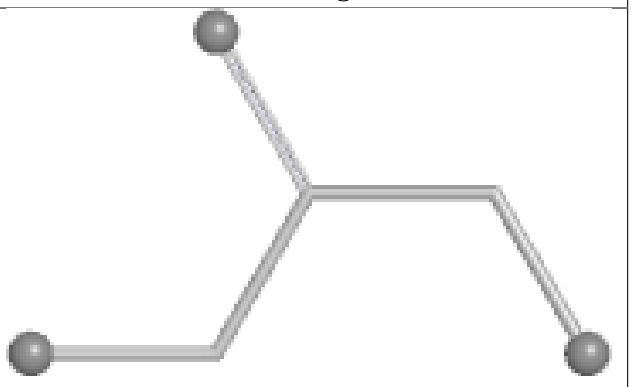
Bond lengths



Bond angles

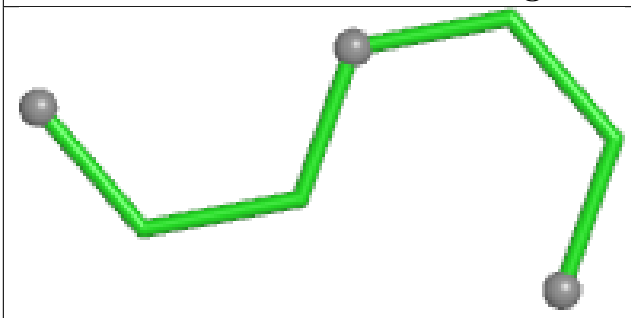


Torsions

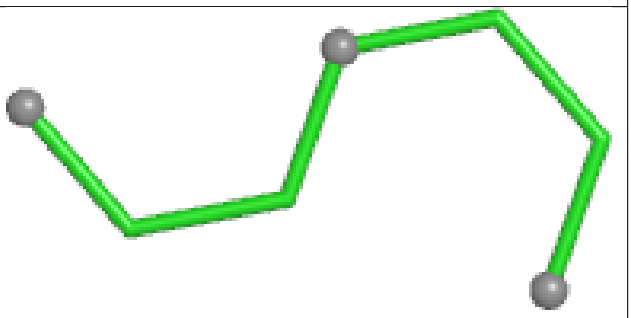


Rings

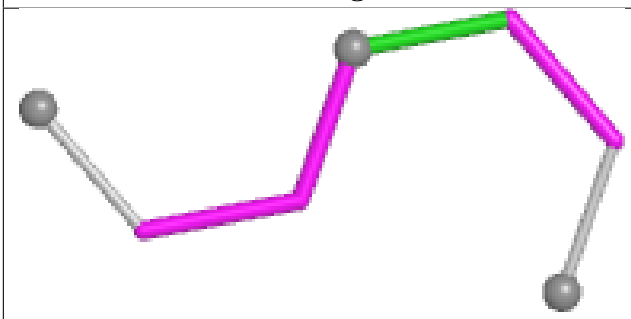
Ligand PEG B 501



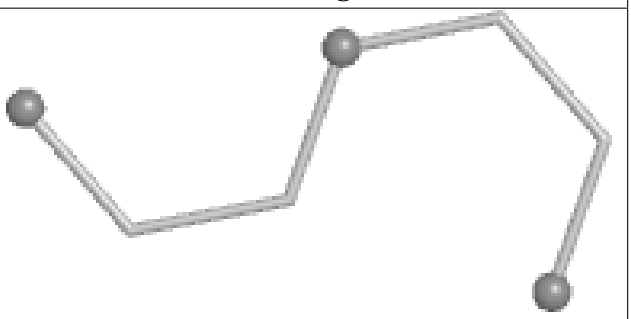
Bond lengths



Bond angles

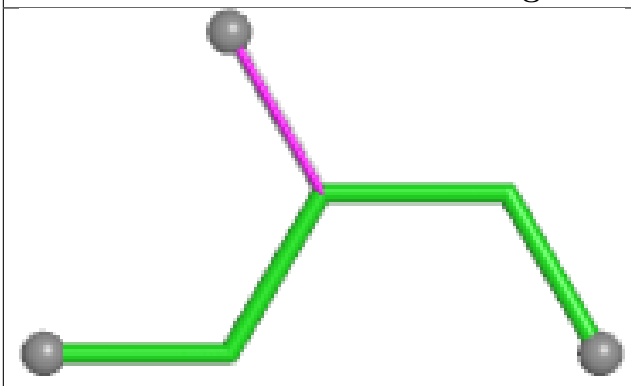


Torsions

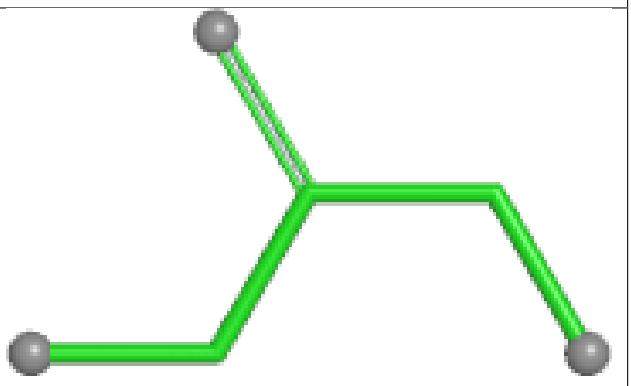


Rings

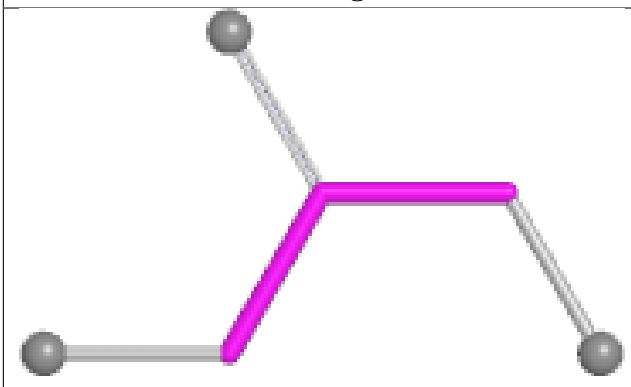
Ligand GOL A 501



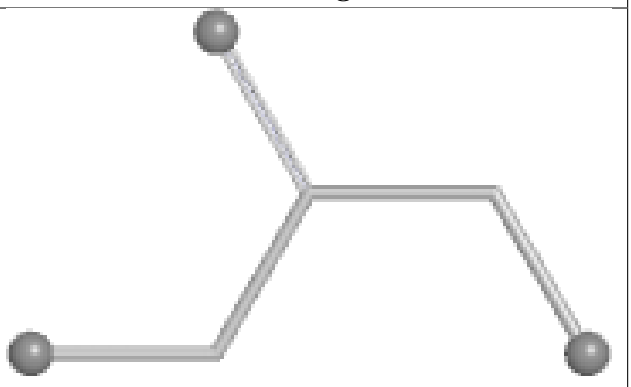
Bond lengths



Bond angles

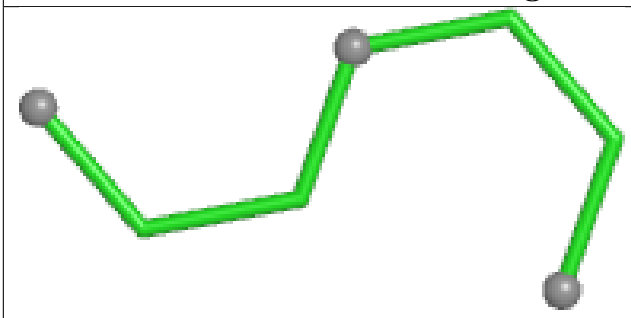


Torsions

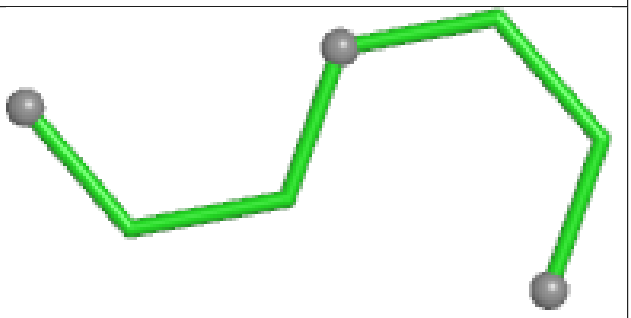


Rings

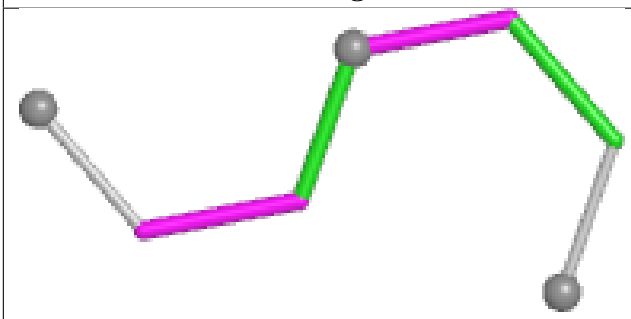
Ligand PEG B 505



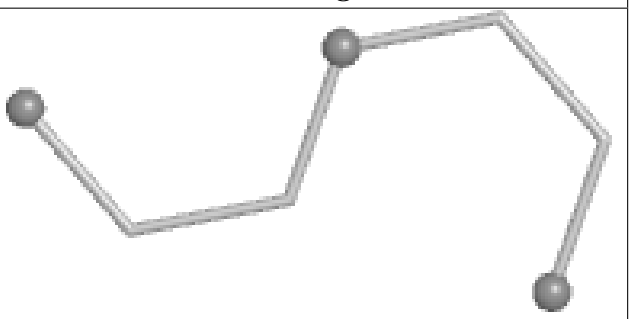
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

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, 95^{th} 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	489/492 (99%)	-0.23	13 (2%)	56	66	23, 37, 59, 140	1 (0%)
1	B	490/492 (99%)	0.00	22 (4%)	38	47	20, 42, 75, 126	1 (0%)
All	All	979/984 (99%)	-0.11	35 (3%)	46	56	20, 39, 69, 140	2 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	SER	5.8
1	B	489	PRO	4.7
1	A	96	MET	4.2
1	B	96	MET	4.2
1	A	450	LEU	3.9
1	B	254	GLY	3.6
1	A	233	LYS	3.2
1	B	450	LEU	3.1
1	A	322	ASP	3.1
1	A	0	SER	3.0
1	B	46	LYS	2.9
1	B	462	ILE	2.9
1	A	345[A]	ARG	2.8
1	B	233	LYS	2.8
1	B	447	LEU	2.7
1	B	453	PRO	2.5
1	A	436	ARG	2.5
1	A	414	TYR	2.4
1	A	126	LEU	2.4
1	B	0	SER	2.4
1	B	286	CYS	2.4
1	B	486	LYS	2.4
1	A	454	PHE	2.3
1	A	254	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	226	ILE	2.2
1	B	246	VAL	2.2
1	A	486	LYS	2.2
1	B	460	SER	2.2
1	B	126	LEU	2.2
1	B	454	PHE	2.1
1	B	449	ASP	2.1
1	B	16	SER	2.0
1	B	17	LYS	2.0
1	B	1	MET	2.0
1	B	436	ARG	2.0

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

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

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} 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.

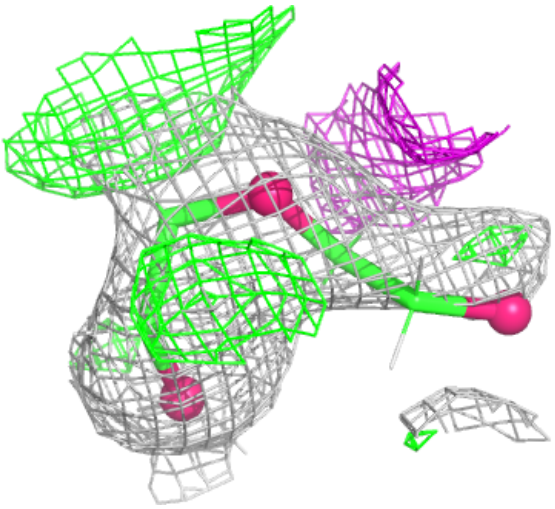
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	PEG	A	504	7/7	0.82	0.22	52,74,115,125	0
4	SO4	B	502	5/5	0.82	0.27	99,105,121,127	0
2	GOL	B	504	6/6	0.83	0.20	45,84,123,123	0
3	PEG	B	505	7/7	0.86	0.21	55,73,93,93	0
3	PEG	A	502	7/7	0.87	0.23	48,75,119,119	0
2	GOL	B	503	6/6	0.87	0.15	73,88,103,105	0
3	PEG	A	503	7/7	0.88	0.21	47,68,121,121	0
3	PEG	B	501	7/7	0.91	0.17	55,69,102,102	0
2	GOL	A	501	6/6	0.92	0.19	37,70,94,94	0

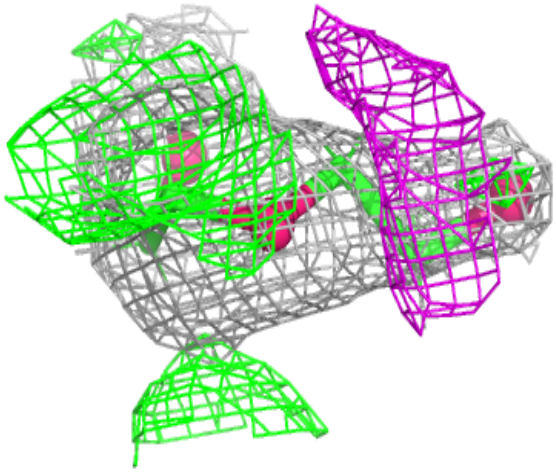
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

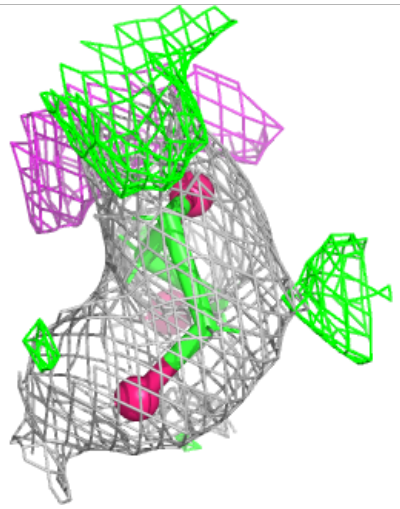
as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PEG A 504:

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

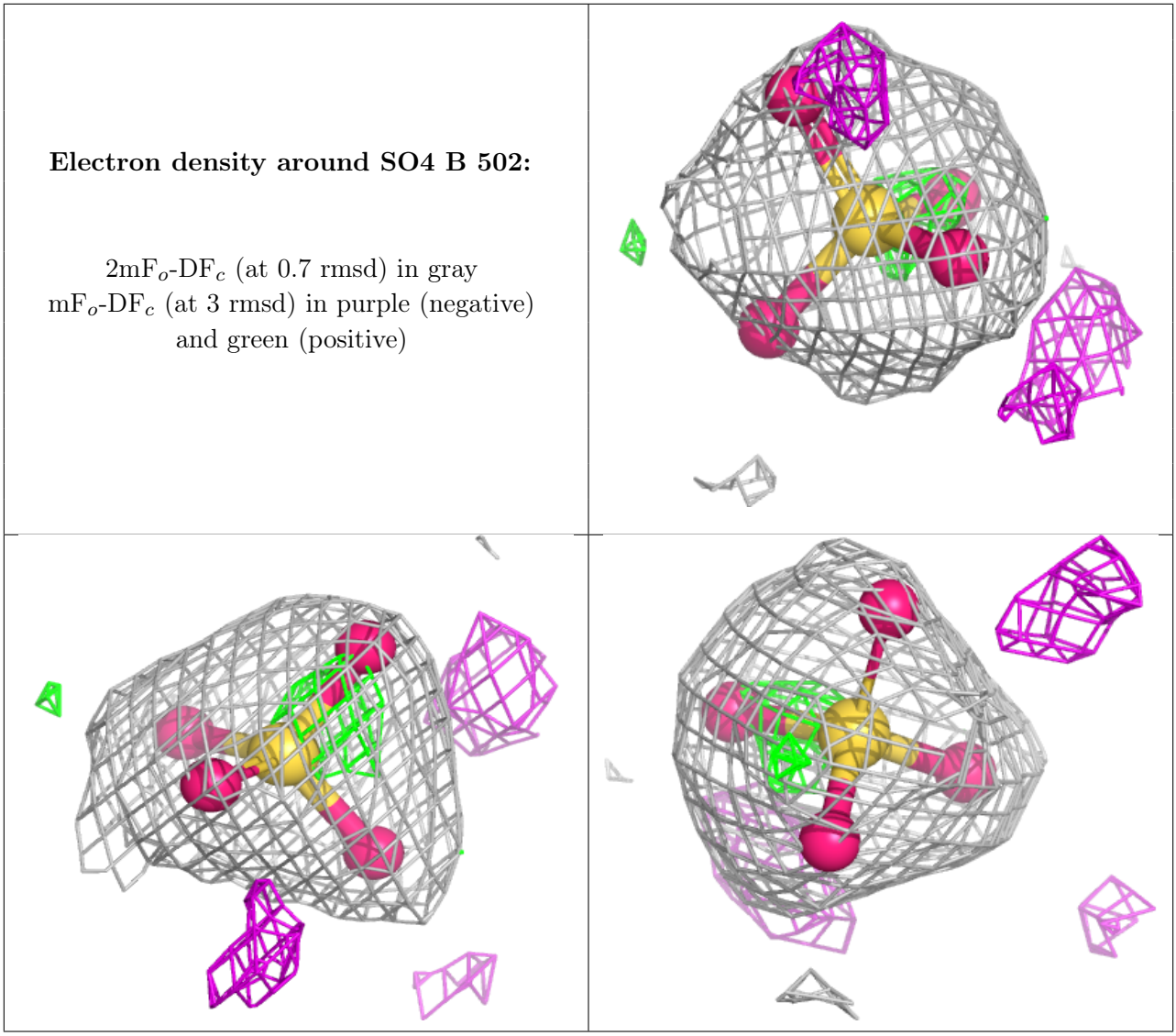






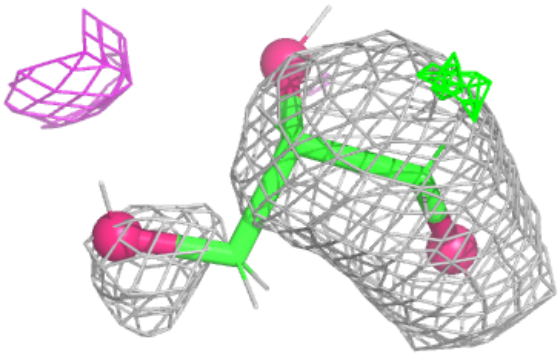
Electron density around SO4 B 502:

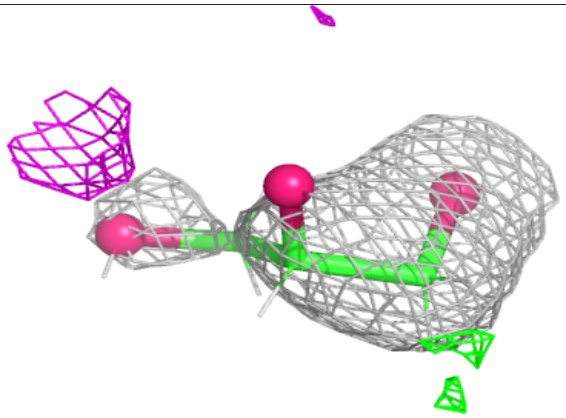
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

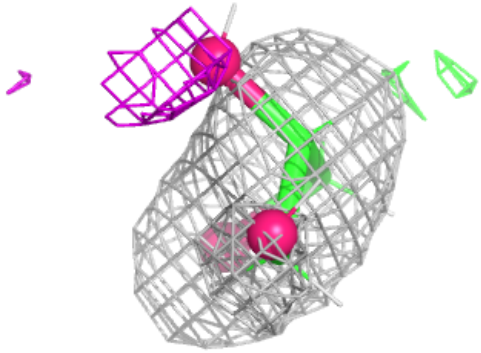


Electron density around GOL B 504:

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

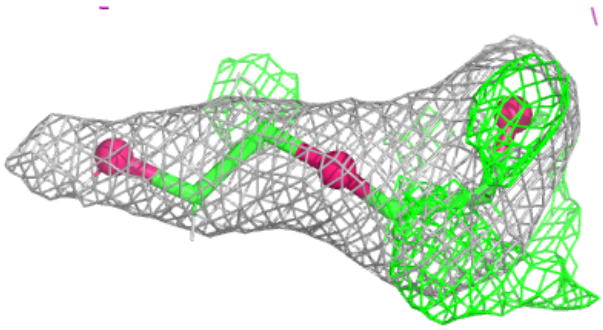


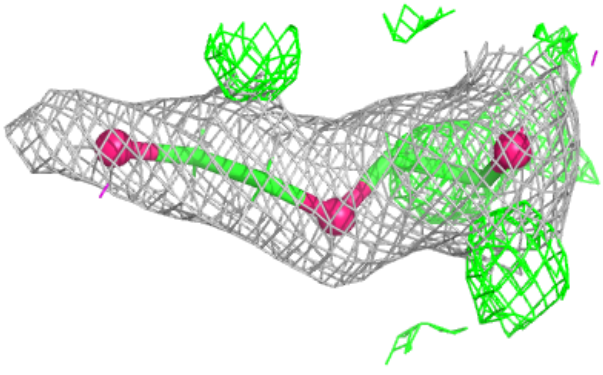


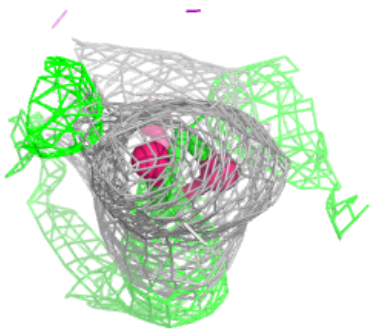


Electron density around PEG B 505:

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

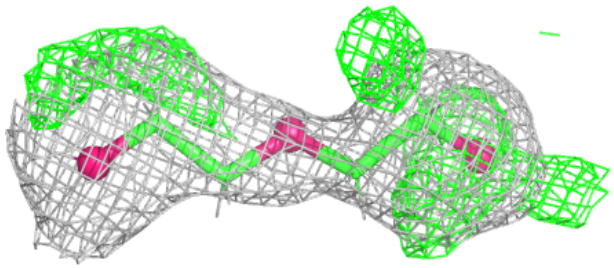


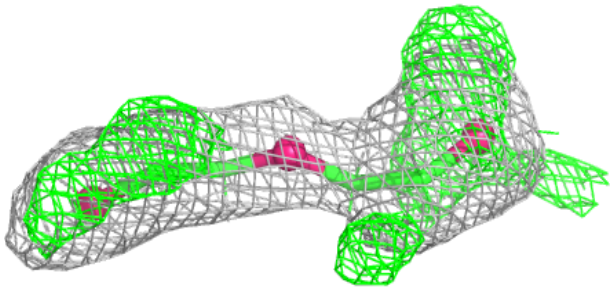


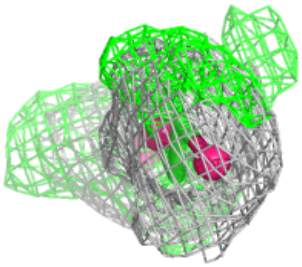


Electron density around PEG A 502:

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

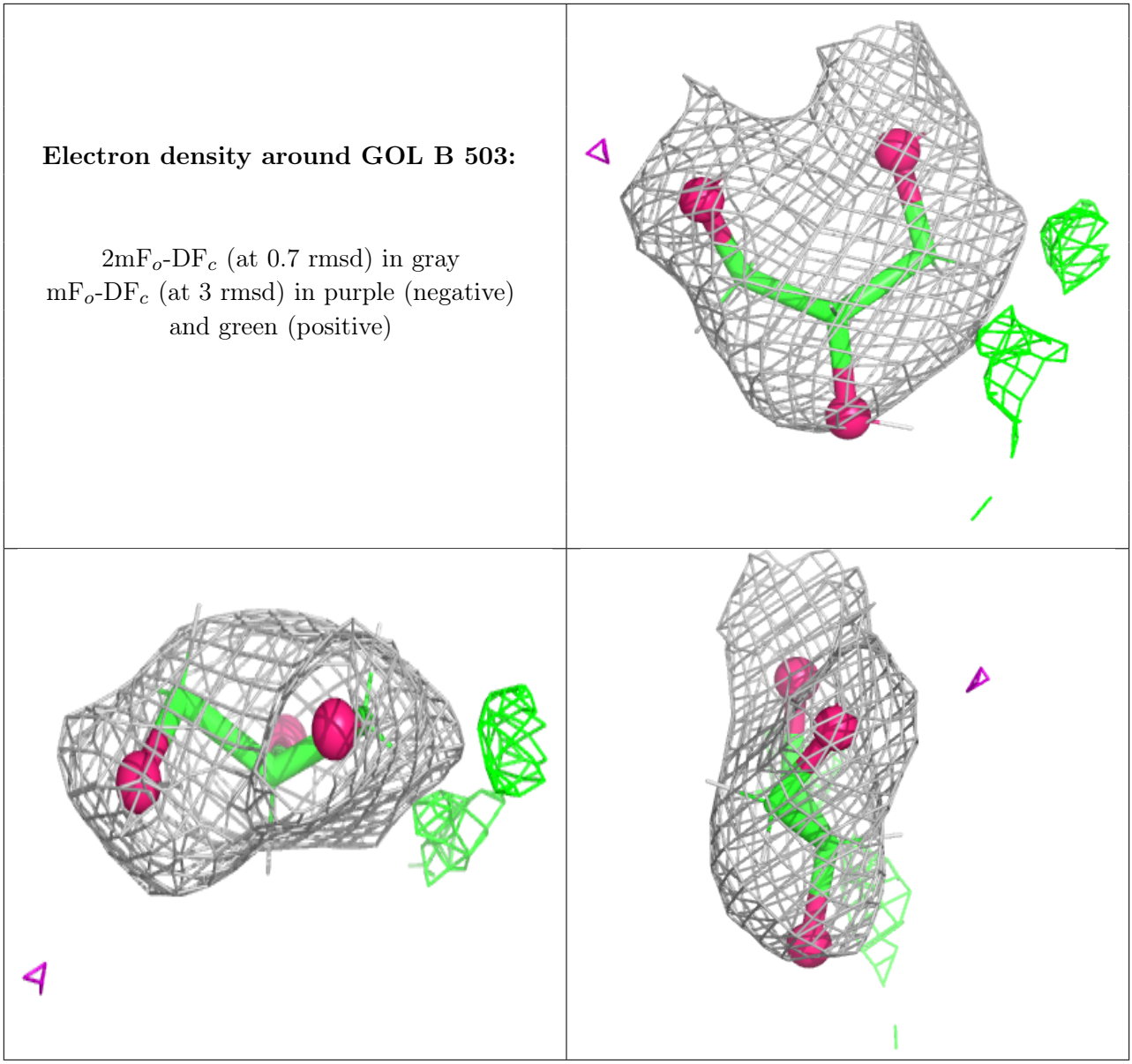






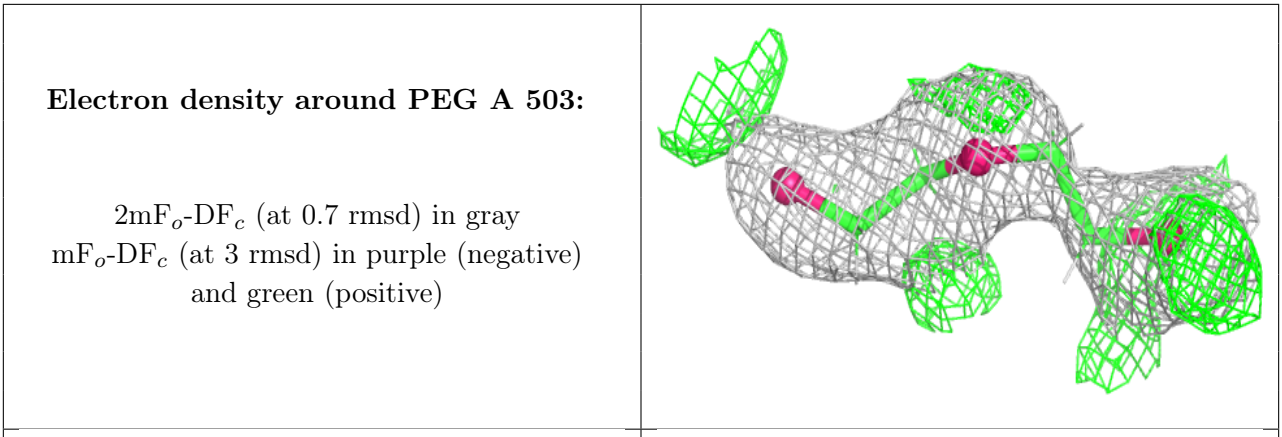
Electron density around GOL B 503:

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



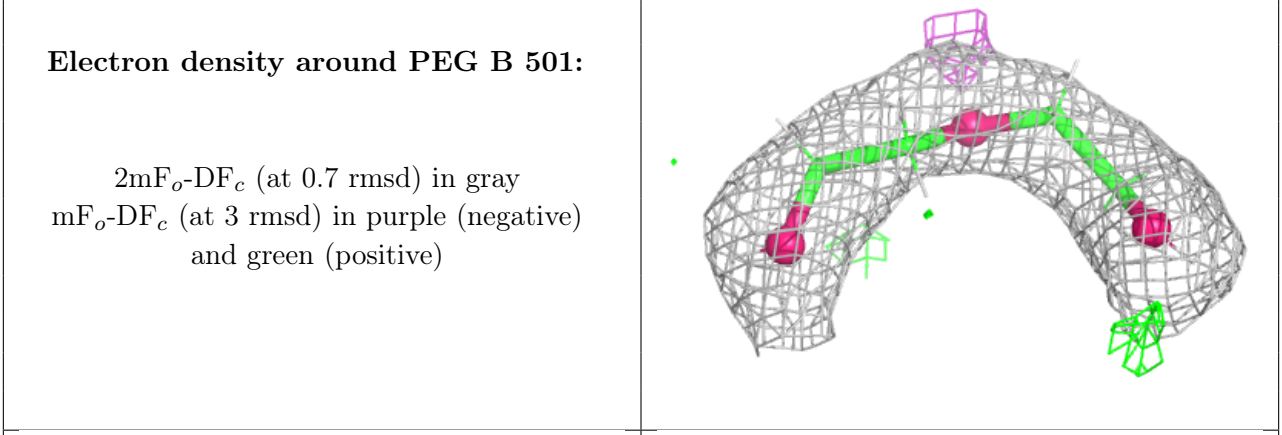
Electron density around PEG A 503:

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



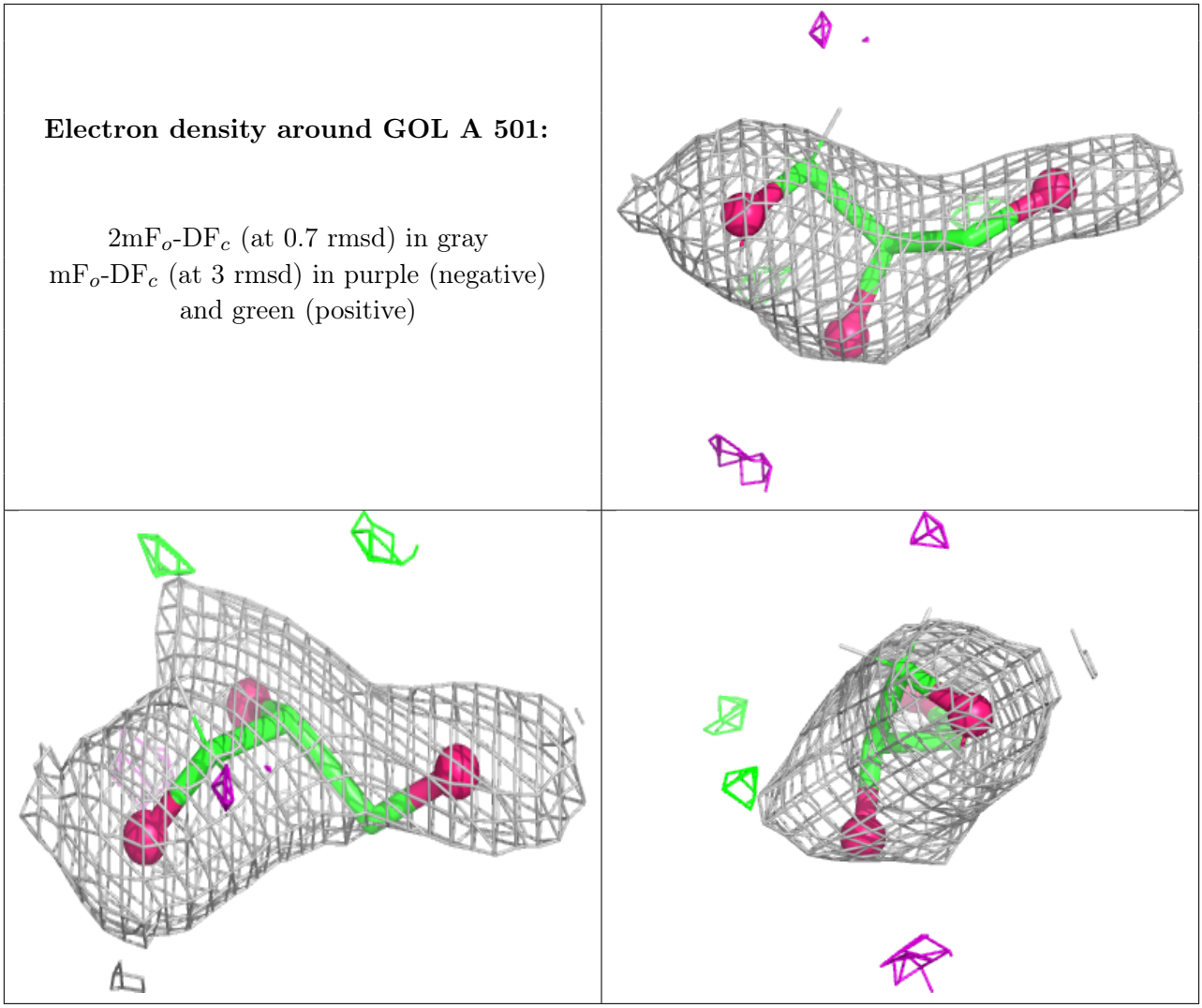
Electron density around PEG B 501:

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



Electron density around GOL A 501:

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



6.5 Other polymers

There are no such residues in this entry.