



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 05:52 AM UTC

PDB ID : 7DL7 / pdb_00007dl7
Title : The wild-type structure of 3,5-DAHDHcca
Authors : Liu, N.; Wu, L.; Zhu, D.M.; Zhou, J.H.
Deposited on : 2020-11-26
Resolution : 2.30 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

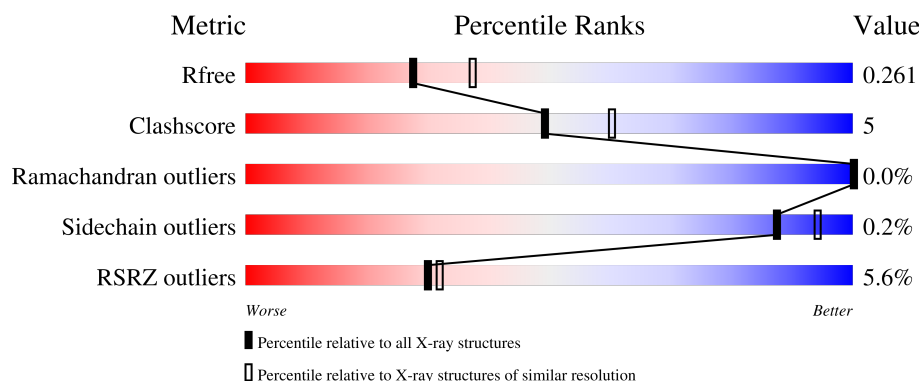
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	358	2% 86% 12% .
1	B	358	3% 84% 14% .
1	C	358	6% 88% 10% .
1	D	358	3% 87% 11% .
1	E	358	14% 81% 16% .

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Mol	Chain	Length	Quality of chain
1	F	358	 4% 85% 13% •
1	G	358	 9% 85% 13% •
1	H	358	 3% 84% 13% ••

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 21840 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 3,5-diaminohexanoate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	350	Total	C	N	O	S	0	1	0
			2672	1676	464	514	18			
1	B	350	Total	C	N	O	S	0	0	0
			2655	1667	459	511	18			
1	C	351	Total	C	N	O	S	0	0	0
			2645	1660	460	507	18			
1	D	350	Total	C	N	O	S	0	1	0
			2655	1667	461	509	18			
1	E	350	Total	C	N	O	S	0	1	0
			2611	1640	453	499	19			
1	F	351	Total	C	N	O	S	0	1	0
			2662	1671	461	511	19			
1	G	351	Total	C	N	O	S	0	2	0
			2654	1663	463	510	18			
1	H	350	Total	C	N	O	S	0	1	0
			2662	1671	460	512	19			

There are 48 discrepancies between the modelled and reference sequences:

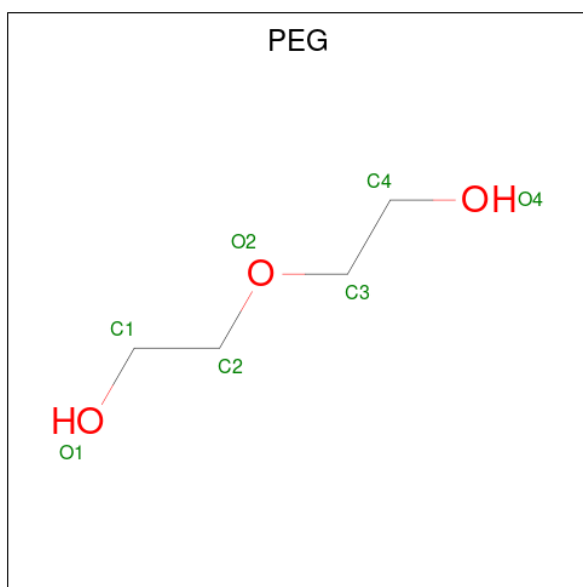
Chain	Residue	Modelled	Actual	Comment	Reference
A	353	HIS	-	expression tag	UNP B0VJ11
A	354	HIS	-	expression tag	UNP B0VJ11
A	355	HIS	-	expression tag	UNP B0VJ11
A	356	HIS	-	expression tag	UNP B0VJ11
A	357	HIS	-	expression tag	UNP B0VJ11
A	358	HIS	-	expression tag	UNP B0VJ11
B	353	HIS	-	expression tag	UNP B0VJ11
B	354	HIS	-	expression tag	UNP B0VJ11
B	355	HIS	-	expression tag	UNP B0VJ11
B	356	HIS	-	expression tag	UNP B0VJ11
B	357	HIS	-	expression tag	UNP B0VJ11
B	358	HIS	-	expression tag	UNP B0VJ11
C	353	HIS	-	expression tag	UNP B0VJ11

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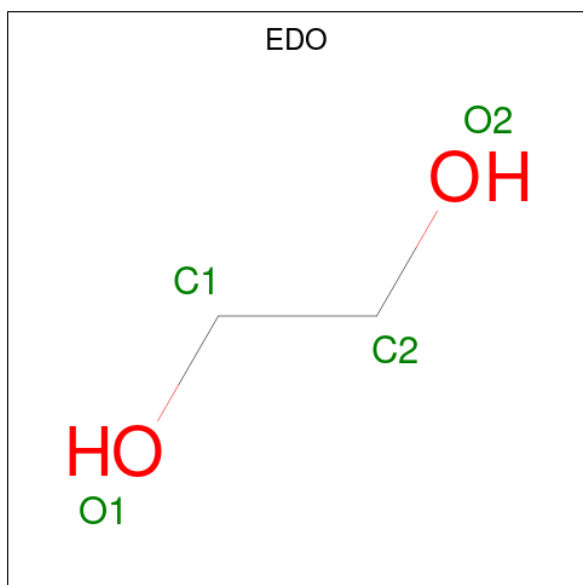
Chain	Residue	Modelled	Actual	Comment	Reference
C	354	HIS	-	expression tag	UNP B0VJ11
C	355	HIS	-	expression tag	UNP B0VJ11
C	356	HIS	-	expression tag	UNP B0VJ11
C	357	HIS	-	expression tag	UNP B0VJ11
C	358	HIS	-	expression tag	UNP B0VJ11
D	353	HIS	-	expression tag	UNP B0VJ11
D	354	HIS	-	expression tag	UNP B0VJ11
D	355	HIS	-	expression tag	UNP B0VJ11
D	356	HIS	-	expression tag	UNP B0VJ11
D	357	HIS	-	expression tag	UNP B0VJ11
D	358	HIS	-	expression tag	UNP B0VJ11
E	353	HIS	-	expression tag	UNP B0VJ11
E	354	HIS	-	expression tag	UNP B0VJ11
E	355	HIS	-	expression tag	UNP B0VJ11
E	356	HIS	-	expression tag	UNP B0VJ11
E	357	HIS	-	expression tag	UNP B0VJ11
E	358	HIS	-	expression tag	UNP B0VJ11
F	353	HIS	-	expression tag	UNP B0VJ11
F	354	HIS	-	expression tag	UNP B0VJ11
F	355	HIS	-	expression tag	UNP B0VJ11
F	356	HIS	-	expression tag	UNP B0VJ11
F	357	HIS	-	expression tag	UNP B0VJ11
F	358	HIS	-	expression tag	UNP B0VJ11
G	353	HIS	-	expression tag	UNP B0VJ11
G	354	HIS	-	expression tag	UNP B0VJ11
G	355	HIS	-	expression tag	UNP B0VJ11
G	356	HIS	-	expression tag	UNP B0VJ11
G	357	HIS	-	expression tag	UNP B0VJ11
G	358	HIS	-	expression tag	UNP B0VJ11
H	353	HIS	-	expression tag	UNP B0VJ11
H	354	HIS	-	expression tag	UNP B0VJ11
H	355	HIS	-	expression tag	UNP B0VJ11
H	356	HIS	-	expression tag	UNP B0VJ11
H	357	HIS	-	expression tag	UNP B0VJ11
H	358	HIS	-	expression tag	UNP B0VJ11

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	C	1	Total	C	O	0	0
			7	4	3		
2	H	1	Total	C	O	0	0
			7	4	3		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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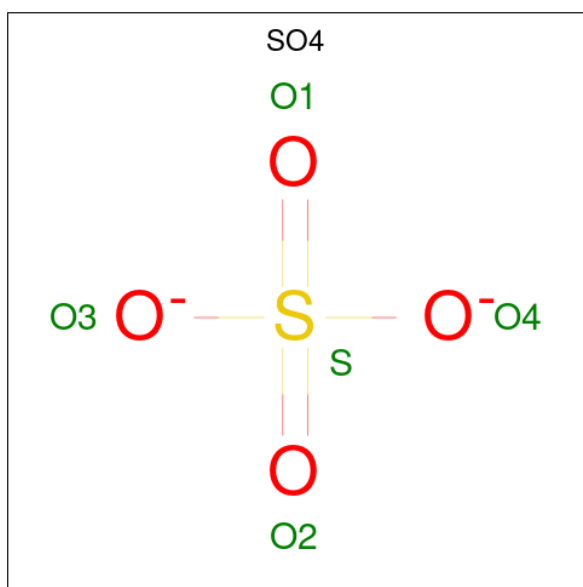
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0
3	H	1	Total C O 4 2 2	0	0

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

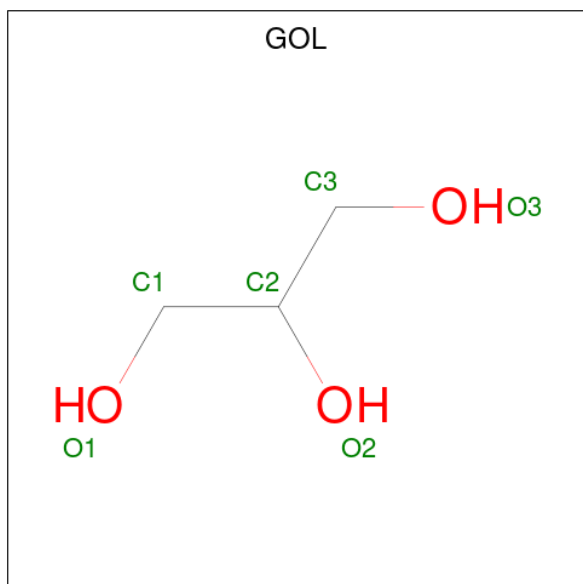
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cl 1 1	0	0
4	C	1	Total Cl 1 1	0	0
4	D	1	Total Cl 1 1	0	0
4	E	1	Total Cl 1 1	0	0
4	F	1	Total Cl 1 1	0	0
4	G	1	Total Cl 1 1	0	0
4	H	2	Total Cl 2 2	0	0

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃) (labeled as "Ligand of Interest" by depositor).

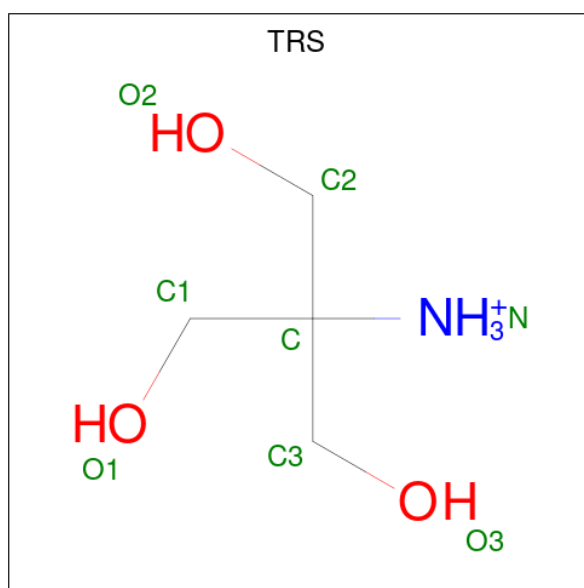


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	1	Total	Ca	0	0
			1	1		
7	G	1	Total	Ca	0	0
			1	1		

- Molecule 8 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	F	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	80	Total	O	0	0
			80	80		
9	B	65	Total	O	0	0
			65	65		
9	C	53	Total	O	0	0
			53	53		
9	D	42	Total	O	0	0
			42	42		
9	E	45	Total	O	0	0
			45	45		

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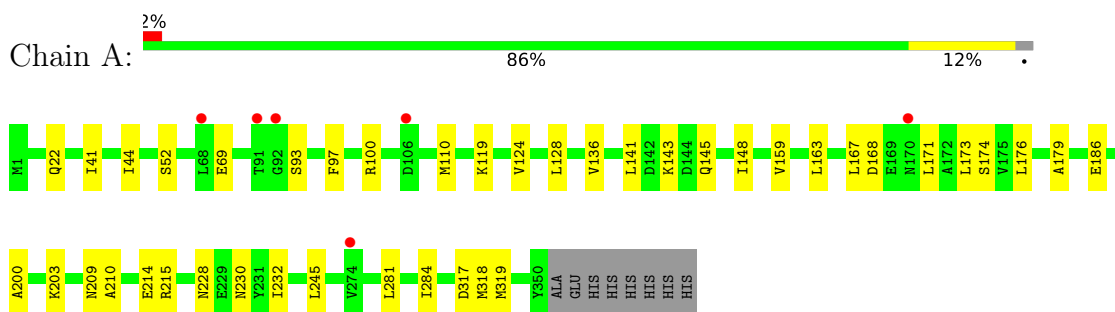
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	F	65	Total 65	O 65	0	0
9	G	43	Total 43	O 43	0	0
9	H	61	Total 61	O 61	0	0

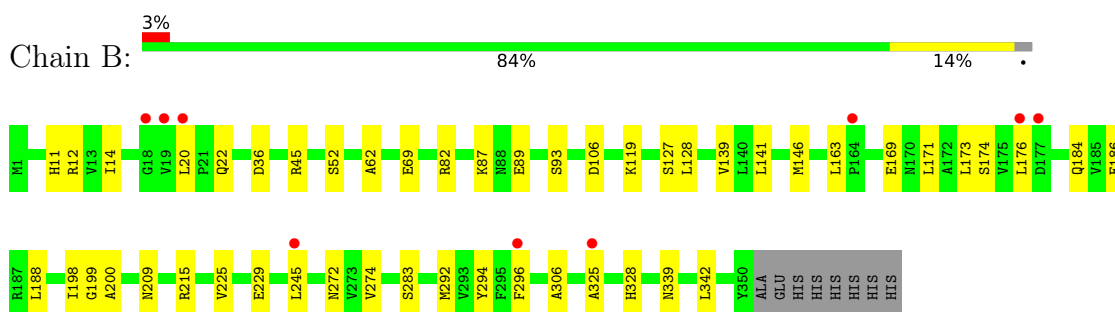
3 Residue-property plots [i](#)

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.

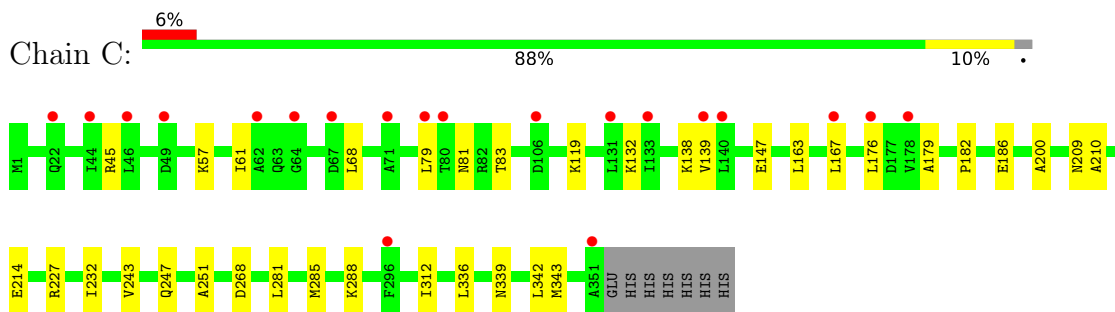
- Molecule 1: 3,5-diaminohexanoate dehydrogenase



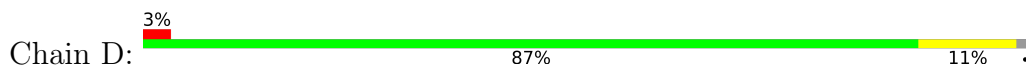
- Molecule 1: 3,5-diaminohexanoate dehydrogenase

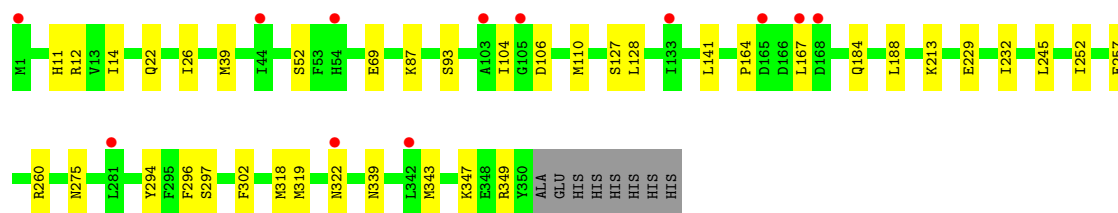


- Molecule 1: 3,5-diaminohexanoate dehydrogenase

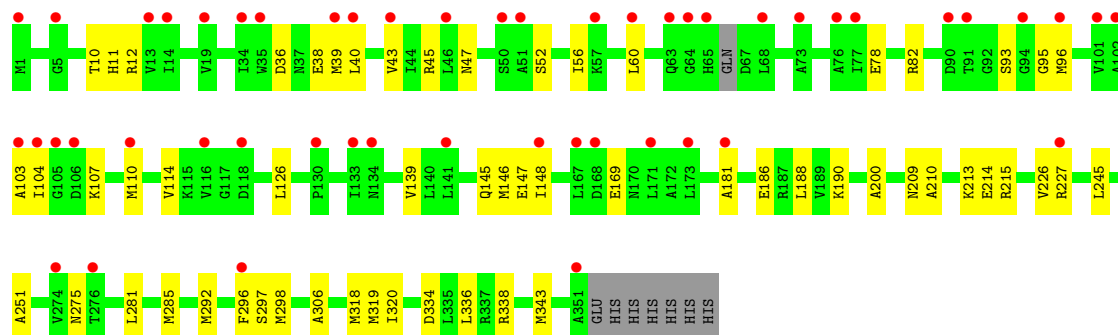
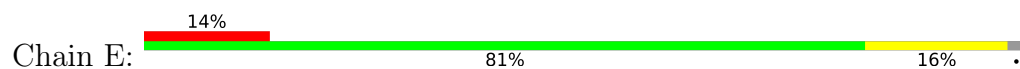


- Molecule 1: 3,5-diaminohexanoate dehydrogenase

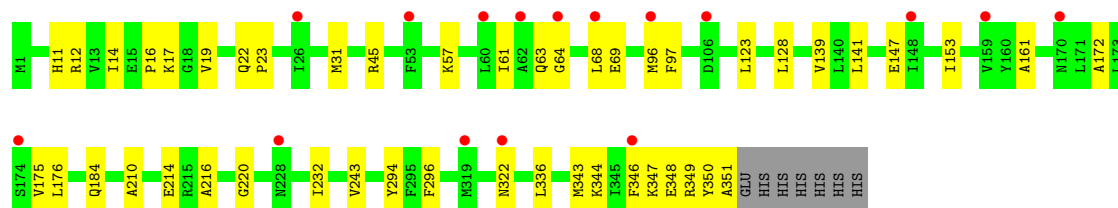
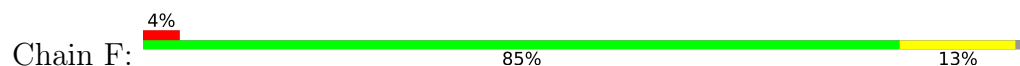




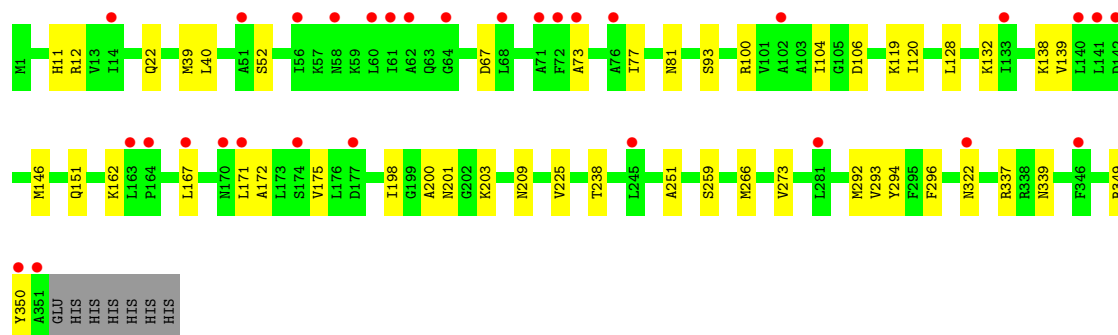
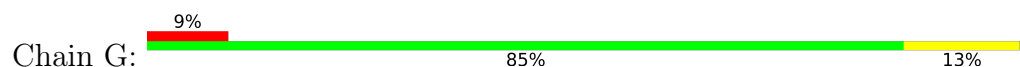
- Molecule 1: 3,5-diaminohexanoate dehydrogenase



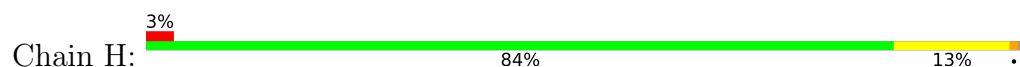
- Molecule 1: 3,5-diaminohexanoate dehydrogenase

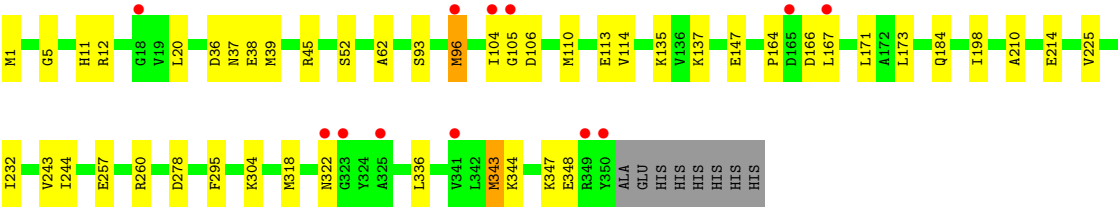


- Molecule 1: 3,5-diaminohexanoate dehydrogenase



- Molecule 1: 3,5-diaminohexanoate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	89.60Å 89.89Å 123.69Å 87.33° 70.97° 80.56°	Depositor
Resolution (Å)	48.98 – 2.30 48.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.3 (48.98-2.30) 97.4 (48.98-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.226 , 0.257 0.229 , 0.261	Depositor DCC
R_{free} test set	7607 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.007 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	21840	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.09	0/2705	0.26	0/3644
1	B	0.08	0/2688	0.25	0/3623
1	C	0.09	0/2678	0.26	0/3612
1	D	0.08	0/2689	0.26	0/3623
1	E	0.12	0/2643	0.32	0/3564
1	F	0.09	0/2695	0.27	0/3632
1	G	0.09	0/2686	0.27	0/3622
1	H	0.09	0/2695	0.27	0/3632
All	All	0.09	0/21479	0.27	0/28952

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	2672	0	2733	24	0
1	B	2655	0	2710	35	0
1	C	2645	0	2686	25	0
1	D	2655	0	2702	27	0
1	E	2611	0	2634	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2662	0	2712	30	0
1	G	2654	0	2682	30	0
1	H	2662	0	2716	32	0
2	A	7	0	10	0	0
2	B	7	0	10	0	0
2	C	7	0	10	0	0
2	H	7	0	10	0	0
3	A	12	0	18	0	0
3	B	20	0	30	0	0
3	C	16	0	24	0	0
3	D	8	0	12	0	0
3	E	4	0	6	0	0
3	F	12	0	18	0	0
3	G	20	0	30	2	0
3	H	16	0	24	0	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	1	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
5	B	5	0	0	0	0
5	F	5	0	0	0	0
6	D	6	0	8	0	0
7	D	1	0	0	0	0
7	G	1	0	0	0	0
8	F	8	0	12	0	0
9	A	80	0	0	0	0
9	B	65	0	0	0	0
9	C	53	0	0	0	0
9	D	42	0	0	1	0
9	E	45	0	0	2	0
9	F	65	0	0	0	0
9	G	43	0	0	0	0
9	H	61	0	0	0	0
All	All	21840	0	21797	231	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 5.

All (231) 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:G:259:SER:HA	1:G:266:MET:HE1	1.59	0.84
1:A:232:ILE:HD11	1:A:245:LEU:HD11	1.59	0.83
1:B:294:TYR:CE2	1:B:296:PHE:CZ	2.71	0.78
1:B:294:TYR:HE2	1:B:296:PHE:CZ	2.03	0.76
1:D:232:ILE:HD11	1:D:245:LEU:HD21	1.67	0.76
1:B:294:TYR:CD2	1:B:296:PHE:CE2	2.73	0.76
1:H:336:LEU:HB3	1:H:343:MET:HG2	1.70	0.71
1:C:57:LYS:HG3	1:C:68:LEU:HD23	1.72	0.70
1:H:110:MET:HE3	1:H:114:VAL:H	1.57	0.69
1:C:227:ARG:HG2	1:C:247:GLN:HA	1.74	0.67
1:E:78:GLU:OE2	1:E:82:ARG:NH1	2.28	0.66
1:C:281:LEU:HG	1:C:285:MET:HE2	1.77	0.66
1:B:294:TYR:CE2	1:B:296:PHE:CE2	2.83	0.66
1:F:343:MET:HE2	1:F:347:LYS:HG3	1.76	0.66
1:D:343:MET:HE3	1:D:347:LYS:HE3	1.77	0.65
1:D:339:ASN:ND2	4:D:403:CL:CL	2.66	0.65
1:A:174:SER:O	1:A:203:LYS:NZ	2.29	0.64
1:G:339:ASN:ND2	3:G:405:EDO:O1	2.31	0.64
1:H:45:ARG:NH1	1:H:147:GLU:OE1	2.31	0.64
1:C:200:ALA:HB1	1:C:209:ASN:HD21	1.64	0.63
1:D:164:PRO:HG2	1:D:167:LEU:HB2	1.82	0.61
1:B:69:GLU:HG2	1:B:141:LEU:HD11	1.82	0.61
1:E:251:ALA:HB1	1:E:285:MET:HE2	1.83	0.61
1:G:167:LEU:HD11	1:G:337:ARG:HE	1.65	0.61
1:G:81:ASN:OD1	1:G:132:LYS:NZ	2.31	0.59
1:B:62:ALA:O	1:H:304:LYS:NZ	2.29	0.59
1:C:285:MET:CG	1:C:312:ILE:HD13	2.33	0.59
1:H:278:ASP:OD1	1:H:304:LYS:NZ	2.36	0.59
1:A:317:ASP:HB3	1:A:319:MET:HE2	1.84	0.58
1:A:281:LEU:HA	1:A:284:ILE:HD12	1.86	0.57
1:H:96[A]:MET:HE1	1:H:173:LEU:HD11	1.85	0.56
1:F:344:LYS:O	1:F:348:GLU:HG3	2.05	0.56
1:E:200:ALA:HB1	1:E:209:ASN:HD21	1.69	0.56
1:D:184:GLN:HE22	1:D:322:ASN:HB3	1.69	0.56
1:G:273:VAL:HG12	1:G:296:PHE:HB3	1.88	0.55
1:E:11:HIS:CD2	1:E:12:ARG:HG3	2.41	0.55
1:C:268:ASP:HA	1:C:288:LYS:HE2	1.88	0.55
1:A:167:LEU:HB3	1:A:171:LEU:HD12	1.88	0.55
1:A:44:ILE:HD11	1:A:100:ARG:HG3	1.89	0.54
1:H:184:GLN:HE22	1:H:322:ASN:HB3	1.72	0.54
1:B:325:ALA:HB3	1:B:328:HIS:HB3	1.89	0.54
1:C:232:ILE:HG12	1:C:243:VAL:HG11	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:ILE:HA	1:B:225:VAL:HB	1.90	0.54
1:F:14:ILE:O	1:F:17:LYS:NZ	2.37	0.54
1:B:306:ALA:O	1:G:322:ASN:ND2	2.40	0.53
1:A:168:ASP:HB3	1:A:171:LEU:HB2	1.89	0.53
1:F:184:GLN:HE22	1:F:322:ASN:HB3	1.72	0.53
1:E:43:VAL:HB	1:E:148:ILE:HD11	1.90	0.53
1:A:200:ALA:HB1	1:A:209:ASN:HD21	1.73	0.53
1:A:52:SER:HB2	1:A:93:SER:HB2	1.91	0.53
1:B:229:GLU:HA	1:B:245:LEU:HD21	1.90	0.53
1:H:344:LYS:O	1:H:348:GLU:HG3	2.07	0.53
1:B:294:TYR:HD2	1:B:296:PHE:CE2	2.26	0.53
1:D:349:ARG:NH1	9:D:503:HOH:O	2.42	0.53
1:E:52:SER:HB2	1:E:93:SER:HB2	1.90	0.53
1:B:36:ASP:OD1	1:B:36:ASP:N	2.41	0.52
1:C:336:LEU:HB3	1:C:343:MET:HE2	1.91	0.52
1:E:188:LEU:HD22	1:E:292:MET:HE3	1.91	0.52
1:G:201:ASN:O	1:G:349:ARG:NH2	2.43	0.52
1:A:176:LEU:HA	1:A:179:ALA:HB2	1.92	0.52
1:C:339:ASN:HB3	1:C:342:LEU:HB2	1.92	0.52
1:E:227:ARG:NH2	9:E:504:HOH:O	2.42	0.52
1:H:52:SER:HB2	1:H:93:SER:HB2	1.92	0.52
1:C:57:LYS:HG3	1:C:68:LEU:CD2	2.40	0.51
1:E:190:LYS:NZ	9:E:502:HOH:O	2.36	0.51
1:B:106:ASP:N	1:B:106:ASP:OD1	2.43	0.51
1:B:20:LEU:HD11	1:G:251:ALA:HB3	1.92	0.51
1:D:52:SER:HB2	1:D:93:SER:HB2	1.92	0.51
1:G:198:ILE:HA	1:G:225:VAL:HB	1.91	0.51
1:H:110:MET:CE	1:H:114:VAL:H	2.23	0.51
1:F:63:GLN:H	1:F:64:GLY:HA2	1.76	0.51
1:F:16:PRO:HG2	1:F:19:VAL:HG21	1.92	0.50
1:A:69:GLU:HG2	1:A:141:LEU:HD21	1.92	0.50
1:D:106:ASP:N	1:D:106:ASP:OD1	2.44	0.50
1:H:135:LYS:HE2	1:H:137:LYS:HG2	1.93	0.50
1:G:259:SER:CA	1:G:266:MET:HE1	2.39	0.50
1:F:63:GLN:N	1:F:64:GLY:HA2	2.27	0.50
1:H:110:MET:HE3	1:H:113:GLU:HA	1.94	0.50
1:F:139:VAL:HG12	1:F:141:LEU:HD12	1.93	0.50
1:C:45:ARG:NH2	1:C:147:GLU:OE2	2.45	0.49
1:G:138:LYS:HD2	1:G:139:VAL:N	2.27	0.49
1:H:257:GLU:OE2	1:H:260:ARG:NH2	2.38	0.49
1:E:334:ASP:O	1:E:338:ARG:HG3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:LEU:O	1:B:174:SER:OG	2.29	0.49
1:C:167:LEU:HD21	1:C:343:MET:HE1	1.93	0.49
1:F:349:ARG:NH2	1:F:350:TYR:OH	2.45	0.49
1:G:100:ARG:NH1	1:G:119:LYS:HB2	2.27	0.49
1:B:52:SER:HB2	1:B:93:SER:HB2	1.95	0.49
1:G:167:LEU:HD11	1:G:337:ARG:NE	2.27	0.49
1:G:106:ASP:OD1	1:G:106:ASP:N	2.45	0.49
1:H:198:ILE:HA	1:H:225:VAL:HB	1.93	0.49
1:A:119:LYS:HG2	1:A:163:LEU:HD12	1.95	0.49
1:B:87:LYS:HB2	1:B:127:SER:HA	1.95	0.49
1:F:31:MET:HE1	1:F:153:ILE:HG12	1.93	0.49
1:G:200:ALA:HB1	1:G:209:ASN:HD21	1.78	0.49
1:B:200:ALA:HB1	1:B:209:ASN:HD21	1.77	0.49
1:G:292:MET:HE3	1:G:293:VAL:H	1.78	0.48
1:B:184:GLN:HB2	1:B:296:PHE:HZ	1.79	0.48
1:C:81:ASN:OD1	1:C:132:LYS:NZ	2.43	0.48
1:F:57:LYS:O	1:F:61:ILE:HG13	2.13	0.48
1:G:52:SER:HB2	1:G:93:SER:HB2	1.96	0.48
1:A:124:VAL:HG13	1:A:159:VAL:H	1.78	0.48
1:D:69:GLU:HG2	1:D:141:LEU:HD21	1.96	0.48
1:G:138:LYS:HD2	1:G:139:VAL:H	1.79	0.48
1:H:171:LEU:HD21	1:H:343:MET:HE1	1.94	0.48
1:D:319:MET:HE2	1:E:319:MET:HE1	1.93	0.48
1:E:56:ILE:O	1:E:60:LEU:HD22	2.14	0.48
1:E:12:ARG:NE	1:E:38:GLU:OE2	2.46	0.48
1:B:82:ARG:NH2	1:B:89:GLU:OE1	2.47	0.48
1:E:47:ASN:ND2	1:E:145:GLN:HG3	2.29	0.47
1:H:39:MET:HE2	1:H:104:ILE:HG12	1.96	0.47
1:G:39:MET:HE2	1:G:104:ILE:HG12	1.96	0.47
1:F:45:ARG:NH1	1:F:147:GLU:OE2	2.48	0.47
1:B:188:LEU:HD22	1:B:292:MET:HE3	1.95	0.47
1:E:336:LEU:HD12	1:E:343:MET:HA	1.95	0.47
1:A:110:MET:HE2	1:A:110:MET:HB2	1.82	0.47
1:A:173:LEU:HD23	1:A:176:LEU:HD12	1.97	0.47
1:H:11:HIS:NE2	1:H:38:GLU:OE2	2.28	0.47
1:G:139:VAL:HG22	1:G:146:MET:HG3	1.96	0.47
1:H:166:ASP:O	1:H:167:LEU:HD23	2.15	0.47
1:E:39:MET:HE2	1:E:104:ILE:HG12	1.97	0.47
1:F:11:HIS:CD2	1:F:12:ARG:HG3	2.49	0.47
1:B:294:TYR:HD2	1:B:296:PHE:CD2	2.33	0.47
1:G:172:ALA:O	1:G:175:VAL:HG12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:226:VAL:HG23	1:E:245:LEU:HD23	1.97	0.46
1:F:294:TYR:CE2	1:F:296:PHE:HB2	2.50	0.46
1:C:251:ALA:HB1	1:C:285:MET:CE	2.45	0.46
1:D:11:HIS:CD2	1:D:12:ARG:HG3	2.50	0.46
1:D:188:LEU:HD21	1:D:319:MET:HE1	1.97	0.46
1:E:306:ALA:HA	1:E:318:MET:HE1	1.97	0.46
1:B:139:VAL:HG22	1:B:146:MET:HG2	1.98	0.46
1:C:79:LEU:O	1:C:83:THR:OG1	2.31	0.46
1:E:210:ALA:O	1:E:214:GLU:HG2	2.15	0.46
1:F:69:GLU:HG2	1:F:141:LEU:CD1	2.46	0.46
1:C:285:MET:HG2	1:C:312:ILE:HD13	1.98	0.46
1:E:45:ARG:HD3	1:E:145:GLN:OE1	2.16	0.46
1:G:40:LEU:HD11	1:G:151:GLN:HB3	1.97	0.46
1:H:11:HIS:CD2	1:H:12:ARG:HG3	2.50	0.46
1:H:164:PRO:HB2	1:H:166:ASP:OD1	2.16	0.45
1:H:244:ILE:HG23	1:H:257:GLU:HG2	1.98	0.45
1:H:295:PHE:HE2	1:H:318:MET:HE2	1.81	0.45
1:A:228:ASN:ND2	1:A:230:ASN:HB2	2.31	0.45
1:B:22:GLN:HB2	1:B:128:LEU:HB3	1.99	0.45
1:B:272:ASN:ND2	1:B:283:SER:OG	2.45	0.45
1:C:61:ILE:HG12	1:C:68:LEU:HD11	1.97	0.45
1:G:120:ILE:HG22	1:G:162:LYS:HA	1.98	0.45
1:E:39:MET:HE1	1:E:110:MET:HE1	1.99	0.45
1:H:232:ILE:HG12	1:H:243:VAL:HG11	1.99	0.45
1:C:285:MET:HE1	1:H:20:LEU:HD11	1.97	0.45
1:H:106:ASP:OD1	1:H:106:ASP:N	2.46	0.45
1:E:186:GLU:HA	1:E:215:ARG:HG3	1.99	0.45
1:F:61:ILE:HG12	1:F:68:LEU:HD21	1.98	0.45
1:H:210:ALA:O	1:H:214:GLU:HG2	2.17	0.44
1:F:175:VAL:HG12	1:F:346:PHE:CE2	2.52	0.44
1:D:39:MET:HE2	1:D:104:ILE:HG12	1.99	0.44
1:F:96[B]:MET:HG2	1:F:97:PHE:N	2.32	0.44
1:D:22:GLN:HB2	1:D:128:LEU:HB3	1.99	0.44
1:A:228:ASN:HD21	1:A:230:ASN:HB2	1.83	0.44
1:F:123:LEU:HD11	1:F:161:ALA:HB2	2.00	0.44
1:F:336:LEU:HD22	1:F:343:MET:HA	2.00	0.44
1:H:184:GLN:NE2	1:H:322:ASN:HB3	2.33	0.44
1:B:173:LEU:HD23	1:B:176:LEU:HD12	1.99	0.44
1:D:87:LYS:HB2	1:D:127:SER:HA	2.00	0.44
1:D:302:PHE:CE2	1:E:320:ILE:HG23	2.53	0.43
1:A:136:VAL:HA	1:A:148:ILE:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:45:ARG:NH1	1:E:169:GLU:OE2	2.51	0.43
1:E:181:ALA:HA	1:E:296:PHE:CZ	2.53	0.43
1:H:343:MET:O	1:H:347:LYS:HG2	2.19	0.43
1:G:22:GLN:HB2	1:G:128:LEU:HB3	2.00	0.43
1:F:172:ALA:O	1:F:176:LEU:HG	2.19	0.43
1:C:176:LEU:HA	1:C:179:ALA:HB2	2.00	0.43
1:E:36:ASP:HB2	1:E:107:LYS:HB2	2.01	0.43
1:E:110:MET:HE2	1:E:114:VAL:H	1.83	0.43
1:D:184:GLN:NE2	1:D:322:ASN:HB3	2.31	0.43
1:E:275:ASN:HA	1:E:297:SER:OG	2.19	0.43
1:G:294:TYR:CE2	1:G:296:PHE:HB2	2.53	0.43
1:H:5:GLY:O	1:H:37:ASN:ND2	2.51	0.43
1:C:119:LYS:HG2	1:C:163:LEU:HD12	2.00	0.43
1:E:213:LYS:HA	1:E:213:LYS:HD2	1.88	0.43
1:B:186:GLU:HA	1:B:215:ARG:HG3	2.01	0.42
1:C:61:ILE:HD11	1:C:68:LEU:HD21	1.99	0.42
1:C:251:ALA:HB1	1:C:285:MET:HE1	2.00	0.42
1:E:281:LEU:O	1:E:285:MET:HG3	2.18	0.42
1:F:22:GLN:HB2	1:F:128:LEU:HB3	2.01	0.42
1:F:232:ILE:HG12	1:F:243:VAL:HG11	2.01	0.42
1:B:11:HIS:CD2	1:B:12:ARG:HG3	2.54	0.42
1:B:45:ARG:NH1	1:B:169:GLU:OE2	2.48	0.42
1:D:14:ILE:HB	1:D:26:ILE:HG22	2.00	0.42
1:D:245:LEU:HD13	1:D:245:LEU:HA	1.89	0.42
1:G:203:LYS:HD2	1:G:350:TYR:CD1	2.55	0.42
1:F:19:VAL:HG12	1:F:23:PRO:HG2	2.02	0.42
1:A:41:ILE:HG21	1:A:97:PHE:HE1	1.84	0.42
1:A:143:LYS:O	1:A:145:GLN:HG3	2.20	0.42
1:A:22:GLN:HB2	1:A:128:LEU:HB3	2.01	0.42
1:A:210:ALA:O	1:A:214:GLU:HG2	2.20	0.42
1:B:294:TYR:HE2	1:B:296:PHE:CE1	2.38	0.42
1:F:210:ALA:O	1:F:214:GLU:HG3	2.19	0.42
1:F:216:ALA:HB1	1:F:220:GLY:HA3	2.02	0.42
1:D:213:LYS:HA	1:D:213:LYS:HD2	1.87	0.42
1:A:284:ILE:HG12	1:A:318:MET:HE3	2.02	0.42
1:E:298:MET:HE3	1:E:298:MET:HB3	1.97	0.42
1:G:11:HIS:CD2	1:G:12:ARG:HG3	2.55	0.42
1:B:119:LYS:NZ	1:B:169:GLU:OE1	2.44	0.41
1:E:40:LEU:HB3	1:E:103:ALA:HB3	2.02	0.41
1:C:182:PRO:O	1:C:186:GLU:HG3	2.20	0.41
1:B:339:ASN:HB3	1:B:342:LEU:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:ILE:HD11	1:H:62:ALA:HB2	2.02	0.41
1:D:294:TYR:CE2	1:D:296:PHE:HB2	2.55	0.41
1:H:36:ASP:OD1	1:H:36:ASP:N	2.54	0.41
1:C:138:LYS:HG3	1:C:139:VAL:N	2.34	0.41
1:D:257:GLU:OE1	1:D:260:ARG:NH2	2.36	0.41
1:F:336:LEU:HD23	1:F:336:LEU:HA	1.92	0.41
1:G:238:THR:C	3:G:405:EDO:H12	2.45	0.41
1:E:95:GLY:O	1:E:126:LEU:N	2.53	0.41
1:E:139:VAL:HG22	1:E:146:MET:HG2	2.03	0.41
1:G:73:ALA:O	1:G:77:ILE:HD12	2.21	0.41
1:A:186:GLU:HA	1:A:215:ARG:HG3	2.03	0.41
1:G:171:LEU:O	1:G:171:LEU:HD22	2.21	0.41
1:C:210:ALA:O	1:C:214:GLU:HG2	2.21	0.41
1:D:252:ILE:HG13	1:E:10:THR:HG21	2.03	0.41
1:F:68:LEU:HB3	1:F:141:LEU:HD21	2.03	0.41
1:B:119:LYS:HE3	1:B:163:LEU:HD13	2.02	0.41
1:D:110:MET:HE2	1:D:110:MET:HB2	2.01	0.41
1:D:318:MET:HE2	1:D:318:MET:HB2	1.90	0.40
1:D:229:GLU:HA	1:D:245:LEU:HD11	2.02	0.40
1:D:275:ASN:HA	1:D:297:SER:HB2	2.02	0.40
1:E:36:ASP:OD2	1:E:36:ASP:N	2.55	0.40
1:H:1:MET:SD	1:H:105:GLY:HA2	2.61	0.40
1:B:199:GLY:HA3	1:B:274:VAL:HG13	2.03	0.40
1:F:347:LYS:O	1:F:351:ALA:HB2	2.21	0.40
1:F:175:VAL:HG23	1:F:176:LEU:HD23	2.02	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	349/358 (98%)	341 (98%)	8 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	348/358 (97%)	344 (99%)	4 (1%)	0	100	100
1	C	349/358 (98%)	339 (97%)	10 (3%)	0	100	100
1	D	349/358 (98%)	342 (98%)	7 (2%)	0	100	100
1	E	347/358 (97%)	337 (97%)	10 (3%)	0	100	100
1	F	350/358 (98%)	342 (98%)	8 (2%)	0	100	100
1	G	351/358 (98%)	342 (97%)	8 (2%)	1 (0%)	36	46
1	H	349/358 (98%)	340 (97%)	9 (3%)	0	100	100
All	All	2792/2864 (98%)	2727 (98%)	64 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	67	ASP

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	290/296 (98%)	290 (100%)	0	100	100
1	B	287/296 (97%)	287 (100%)	0	100	100
1	C	282/296 (95%)	282 (100%)	0	100	100
1	D	286/296 (97%)	286 (100%)	0	100	100
1	E	276/296 (93%)	273 (99%)	3 (1%)	65	81
1	F	286/296 (97%)	286 (100%)	0	100	100
1	G	282/296 (95%)	282 (100%)	0	100	100
1	H	288/296 (97%)	285 (99%)	3 (1%)	68	82
All	All	2277/2368 (96%)	2271 (100%)	6 (0%)	87	93

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	96[A]	MET
1	E	96[B]	MET
1	E	147	GLU
1	H	96[A]	MET
1	H	96[B]	MET
1	H	343	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	184	GLN
1	A	247	GLN
1	B	228	ASN
1	C	151	GLN
1	C	201	ASN
1	D	247	GLN
1	E	275	ASN
1	E	327	HIS
1	F	327	HIS
1	G	145	GLN
1	G	201	ASN
1	G	209	ASN
1	G	275	ASN
1	G	322	ASN
1	G	327	HIS
1	H	327	HIS

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

Of 45 ligands modelled in this entry, 10 are monoatomic - leaving 35 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	$\# Z > 2$
3	EDO	C	404	-	3,3,3	0.43	0	2,2,2	0.38	0
5	SO4	B	407	-	4,4,4	0.24	0	6,6,6	0.07	0
3	EDO	D	402	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	F	401	-	3,3,3	0.43	0	2,2,2	0.35	0
2	PEG	B	401	-	6,6,6	0.45	0	5,5,5	0.29	0
3	EDO	F	402	-	3,3,3	0.43	0	2,2,2	0.37	0
3	EDO	A	404	-	3,3,3	0.43	0	2,2,2	0.37	0
3	EDO	G	401	-	3,3,3	0.42	0	2,2,2	0.38	0
3	EDO	G	402	-	3,3,3	0.43	0	2,2,2	0.36	0
3	EDO	B	406	-	3,3,3	0.42	0	2,2,2	0.40	0
3	EDO	B	403	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	D	401	-	3,3,3	0.42	0	2,2,2	0.38	0
3	EDO	A	403	-	3,3,3	0.43	0	2,2,2	0.37	0
3	EDO	B	402	-	3,3,3	0.42	0	2,2,2	0.39	0
2	PEG	H	401	-	6,6,6	0.44	0	5,5,5	0.29	0
3	EDO	H	404	-	3,3,3	0.43	0	2,2,2	0.39	0
3	EDO	G	403	-	3,3,3	0.43	0	2,2,2	0.37	0
3	EDO	G	405	-	3,3,3	0.42	0	2,2,2	0.37	0
5	SO4	F	406	-	4,4,4	0.24	0	6,6,6	0.08	0
3	EDO	C	403	-	3,3,3	0.42	0	2,2,2	0.40	0
3	EDO	F	403	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	H	402	-	3,3,3	0.43	0	2,2,2	0.35	0
3	EDO	B	405	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	A	402	-	3,3,3	0.42	0	2,2,2	0.39	0
3	EDO	C	402	-	3,3,3	0.42	0	2,2,2	0.39	0
3	EDO	H	403	-	3,3,3	0.43	0	2,2,2	0.38	0
2	PEG	A	401	-	6,6,6	0.43	0	5,5,5	0.33	0
3	EDO	E	401	-	3,3,3	0.43	0	2,2,2	0.36	0
2	PEG	C	401	-	6,6,6	0.43	0	5,5,5	0.31	0
3	EDO	G	404	-	3,3,3	0.43	0	2,2,2	0.36	0
3	EDO	H	405	-	3,3,3	0.43	0	2,2,2	0.38	0
3	EDO	B	404	-	3,3,3	0.42	0	2,2,2	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	C	405	-	3,3,3	0.43	0	2,2,2	0.38	0
6	GOL	D	404	-	5,5,5	0.36	0	5,5,5	0.42	0
8	TRS	F	404	-	7,7,7	0.33	0	9,9,9	0.31	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	C	404	-	-	0/1/1/1	-
3	EDO	D	402	-	-	0/1/1/1	-
3	EDO	F	401	-	-	0/1/1/1	-
2	PEG	B	401	-	-	0/4/4/4	-
3	EDO	F	402	-	-	0/1/1/1	-
3	EDO	A	404	-	-	0/1/1/1	-
3	EDO	G	401	-	-	0/1/1/1	-
3	EDO	G	402	-	-	0/1/1/1	-
3	EDO	B	406	-	-	0/1/1/1	-
3	EDO	B	403	-	-	0/1/1/1	-
3	EDO	D	401	-	-	0/1/1/1	-
3	EDO	A	403	-	-	0/1/1/1	-
3	EDO	B	402	-	-	0/1/1/1	-
2	PEG	H	401	-	-	0/4/4/4	-
3	EDO	H	404	-	-	0/1/1/1	-
3	EDO	G	403	-	-	0/1/1/1	-
3	EDO	G	405	-	-	0/1/1/1	-
3	EDO	C	403	-	-	0/1/1/1	-
3	EDO	F	403	-	-	0/1/1/1	-
3	EDO	H	402	-	-	0/1/1/1	-
3	EDO	B	405	-	-	0/1/1/1	-
3	EDO	A	402	-	-	0/1/1/1	-
3	EDO	C	402	-	-	0/1/1/1	-
3	EDO	H	403	-	-	0/1/1/1	-
2	PEG	A	401	-	-	0/4/4/4	-
3	EDO	E	401	-	-	0/1/1/1	-
2	PEG	C	401	-	-	0/4/4/4	-
3	EDO	G	404	-	-	0/1/1/1	-
3	EDO	H	405	-	-	1/1/1/1	-
3	EDO	B	404	-	-	1/1/1/1	-
3	EDO	C	405	-	-	0/1/1/1	-
6	GOL	D	404	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	TRS	F	404	-	-	5/9/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

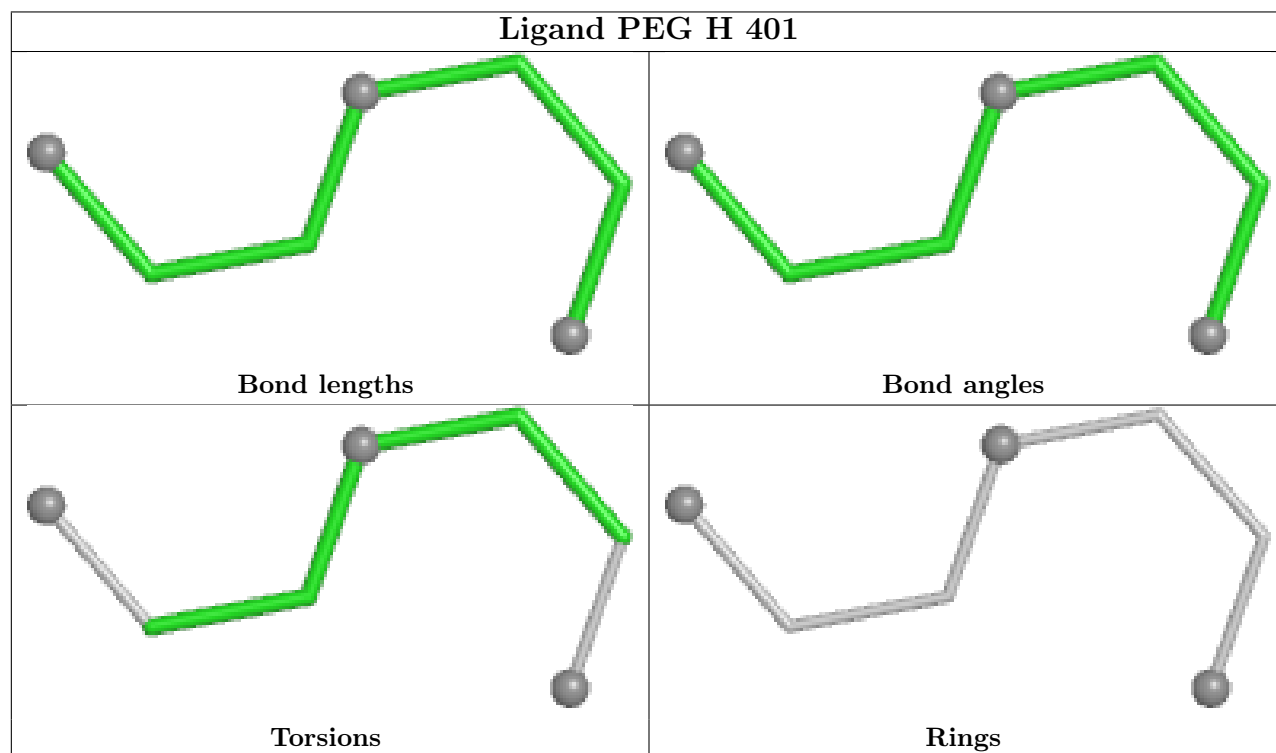
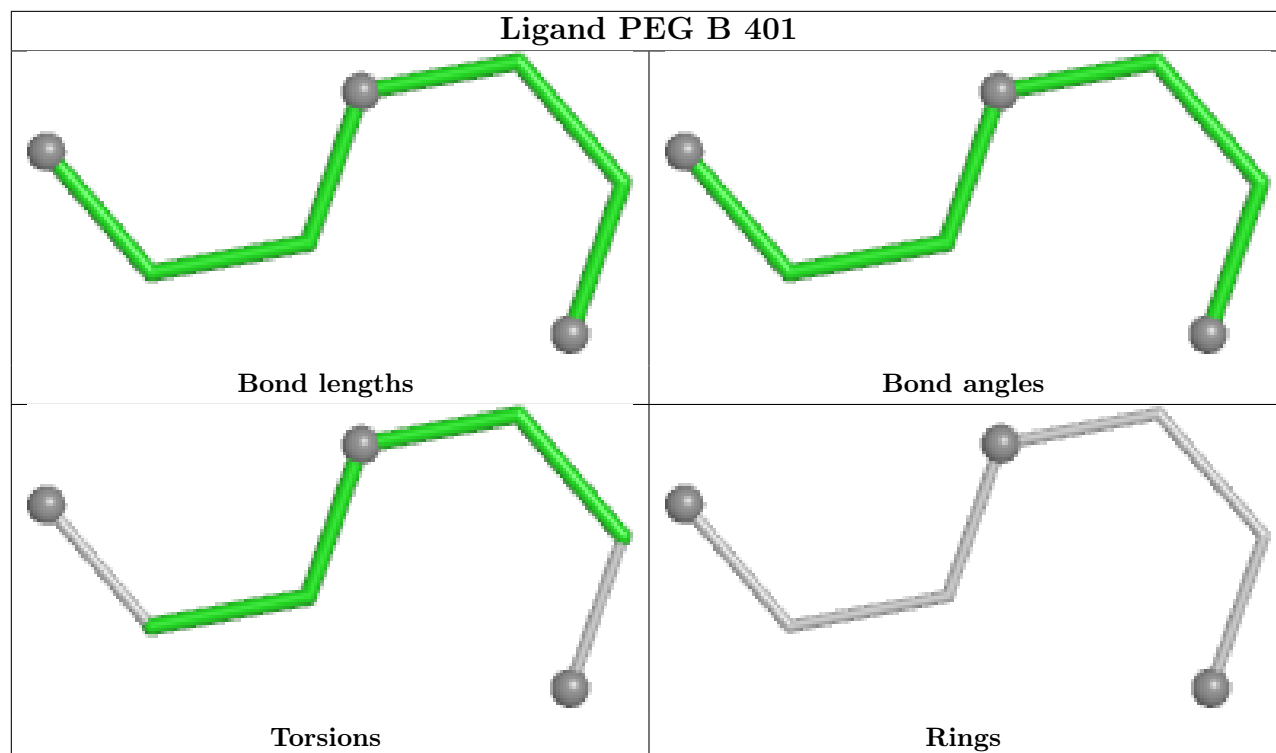
Mol	Chain	Res	Type	Atoms
8	F	404	TRS	C3-C-C2-O2
8	F	404	TRS	N-C-C2-O2
8	F	404	TRS	C1-C-C2-O2
8	F	404	TRS	C2-C-C1-O1
3	B	404	EDO	O1-C1-C2-O2
8	F	404	TRS	N-C-C1-O1
3	H	405	EDO	O1-C1-C2-O2

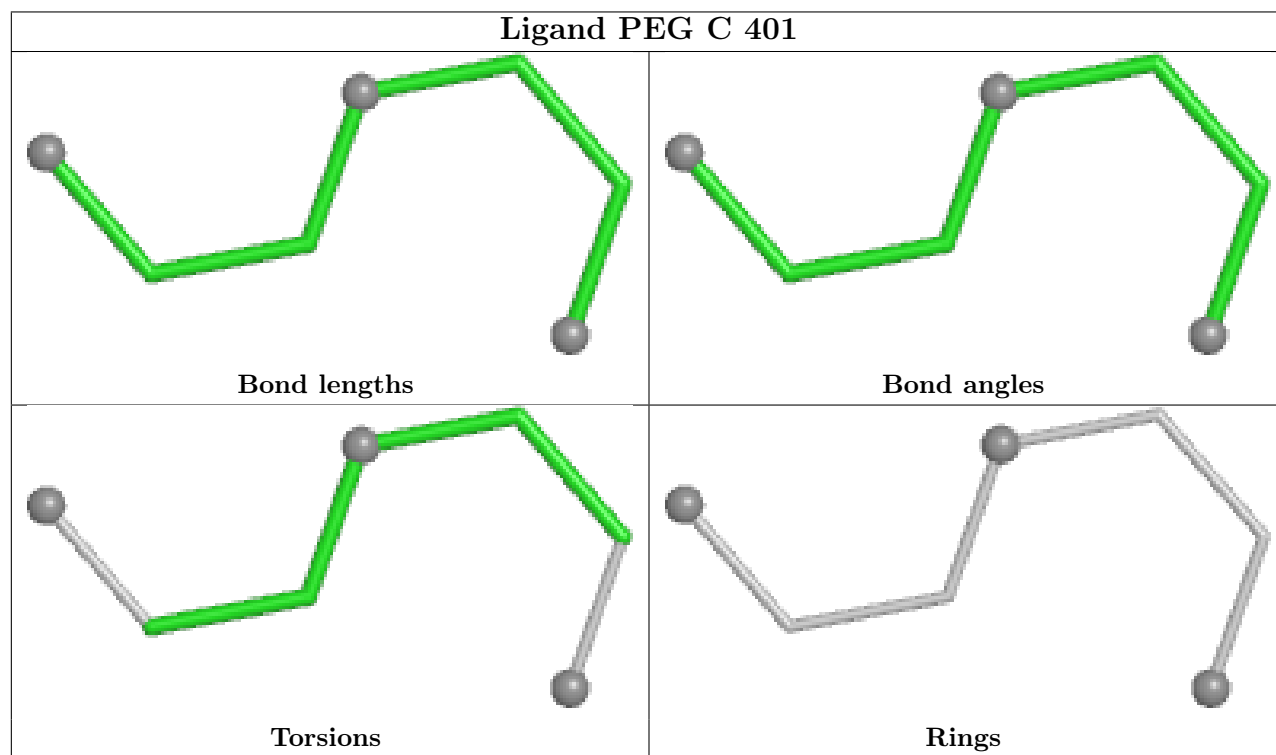
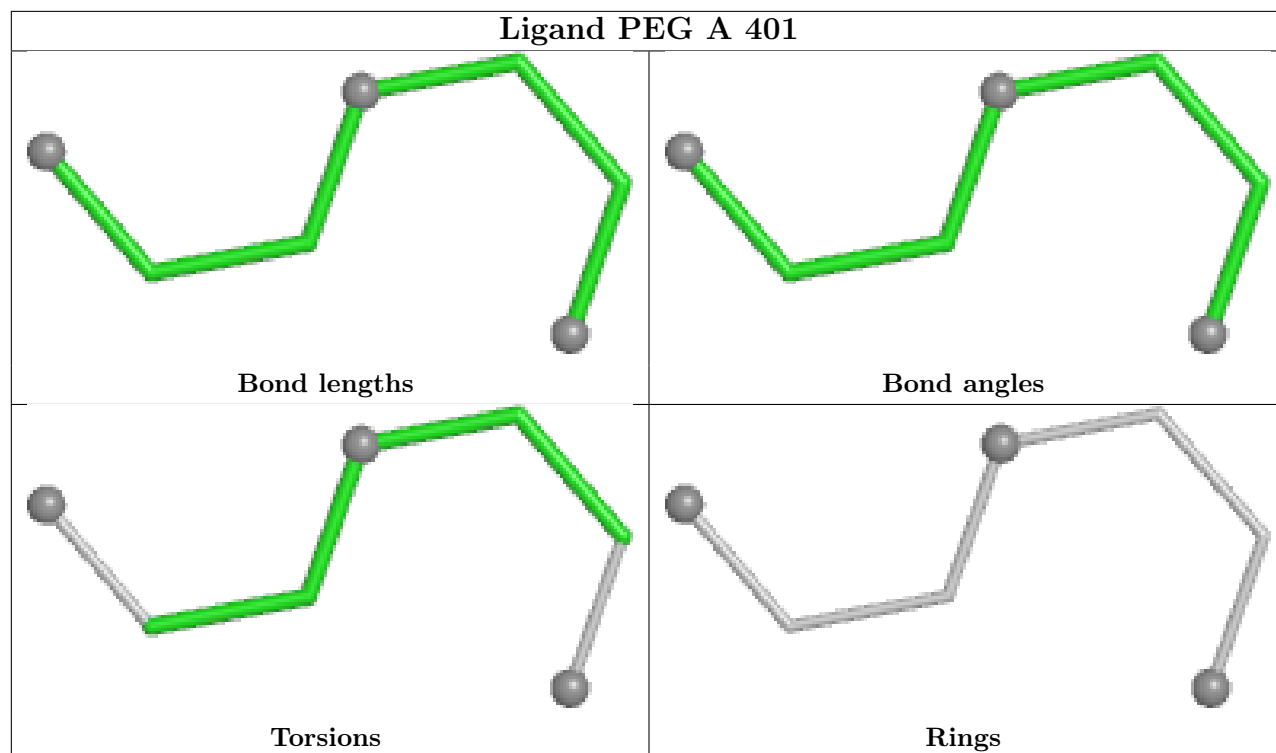
There are no ring outliers.

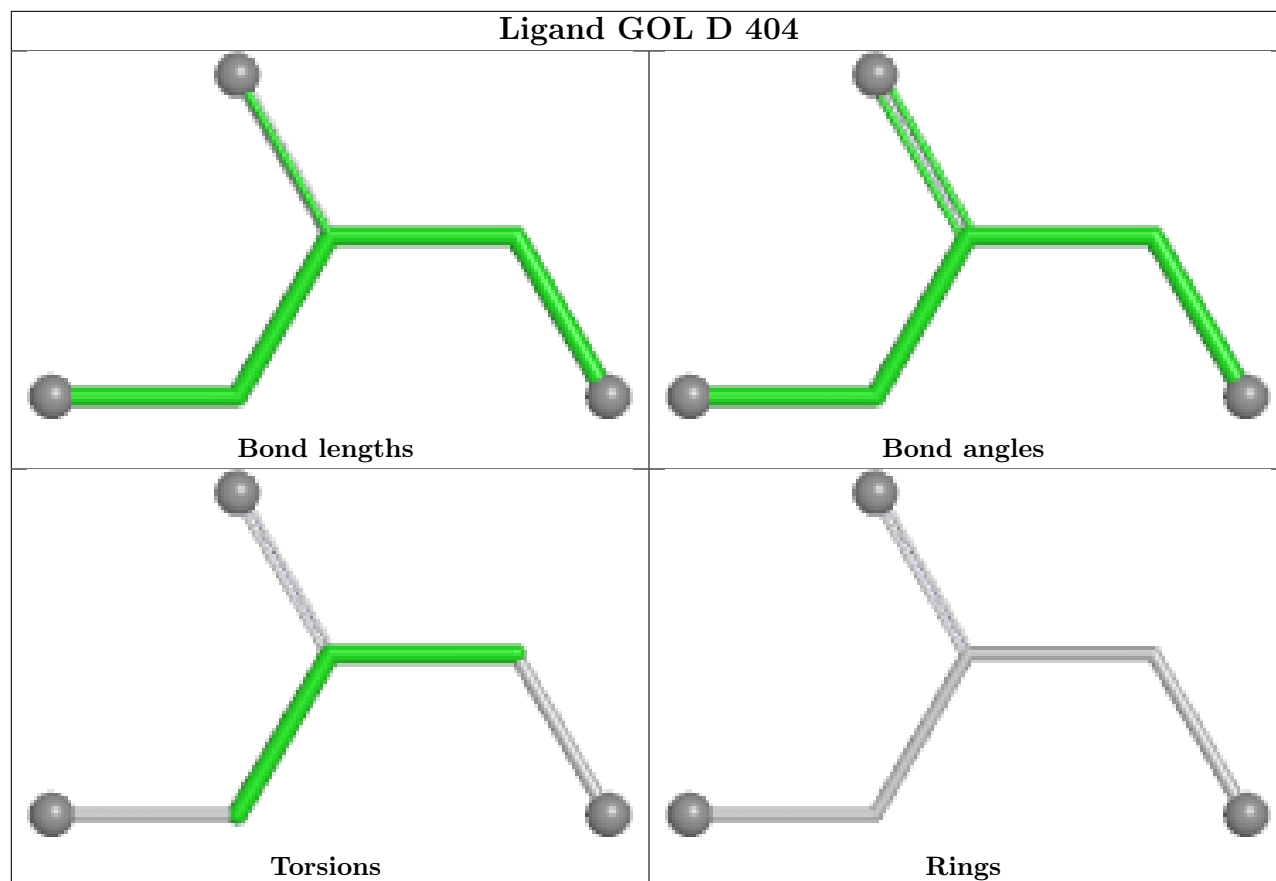
1 monomer is involved in 2 short contacts:

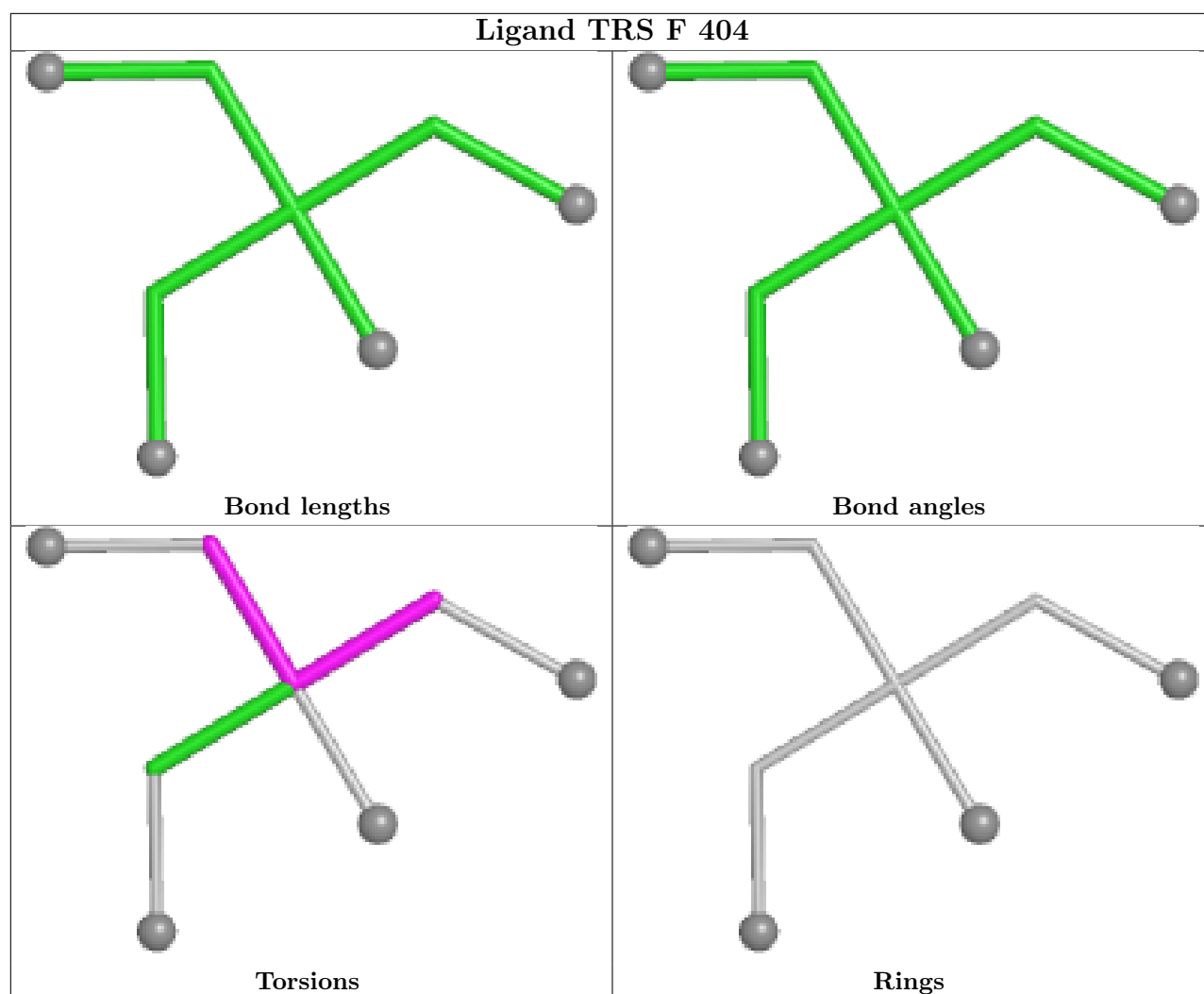
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	405	EDO	2	0

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 [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	350/358 (97%)	0.43	6 (1%) 69 70	18, 41, 54, 73	1 (0%)
1	B	350/358 (97%)	0.51	9 (2%) 57 59	30, 44, 58, 73	0
1	C	351/358 (98%)	0.71	20 (5%) 29 31	30, 47, 66, 85	0
1	D	350/358 (97%)	0.72	12 (3%) 48 50	26, 49, 64, 77	1 (0%)
1	E	350/358 (97%)	1.01	50 (14%) 6 7	28, 52, 72, 81	1 (0%)
1	F	351/358 (98%)	0.65	16 (4%) 37 39	24, 43, 63, 77	1 (0%)
1	G	351/358 (98%)	0.87	31 (8%) 15 17	19, 47, 75, 84	2 (0%)
1	H	350/358 (97%)	0.50	12 (3%) 48 50	20, 41, 58, 73	1 (0%)
All	All	2803/2864 (97%)	0.68	156 (5%) 30 32	18, 45, 67, 85	7 (0%)

All (156) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	104	ILE	4.2
1	A	92	GLY	4.1
1	G	351	ALA	3.9
1	E	1	MET	3.8
1	E	116	VAL	3.8
1	B	296	PHE	3.7
1	E	13	VAL	3.5
1	C	351	ALA	3.5
1	E	102	ALA	3.5
1	F	96[A]	MET	3.4
1	E	91	THR	3.4
1	E	43	VAL	3.4
1	C	64	GLY	3.4
1	F	64	GLY	3.4
1	G	142	ASP	3.4
1	F	228	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	19	VAL	3.3
1	B	325	ALA	3.3
1	G	171	LEU	3.3
1	E	110	MET	3.2
1	E	65	HIS	3.2
1	G	61	ILE	3.2
1	F	159	VAL	3.2
1	B	18	GLY	3.2
1	E	103	ALA	3.2
1	H	96[A]	MET	3.2
1	E	19	VAL	3.2
1	E	351	ALA	3.1
1	G	164	PRO	3.1
1	G	60	LEU	3.1
1	G	73	ALA	3.1
1	E	90	ASP	3.0
1	G	170	ASN	3.0
1	H	105	GLY	3.0
1	H	322	ASN	3.0
1	D	54[A]	HIS	3.0
1	G	174	SER	3.0
1	E	64	GLY	3.0
1	E	40	LEU	2.9
1	G	56	ILE	2.9
1	C	67	ASP	2.9
1	E	96[A]	MET	2.8
1	G	102	ALA	2.8
1	H	167	LEU	2.8
1	G	72	PHE	2.8
1	D	168	ASP	2.7
1	E	34	ILE	2.7
1	E	105	GLY	2.7
1	E	167	LEU	2.7
1	H	165	ASP	2.7
1	B	176	LEU	2.7
1	D	167	LEU	2.7
1	G	141	LEU	2.7
1	G	167	LEU	2.7
1	F	322	ASN	2.7
1	G	64	GLY	2.7
1	D	281	LEU	2.7
1	E	171	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	G	62	ALA	2.6
1	E	60	LEU	2.6
1	F	68	LEU	2.6
1	G	245	LEU	2.6
1	B	164	PRO	2.6
1	C	133	ILE	2.6
1	E	173	LEU	2.6
1	E	76	ALA	2.6
1	E	276	THR	2.6
1	E	133	ILE	2.6
1	H	350	TYR	2.6
1	C	71	ALA	2.5
1	A	274	VAL	2.5
1	C	79	LEU	2.5
1	G	68	LEU	2.5
1	D	165	ASP	2.5
1	E	35	TRP	2.5
1	G	350	TYR	2.5
1	C	176	LEU	2.5
1	D	133	ILE	2.4
1	G	51	ALA	2.4
1	B	245	LEU	2.4
1	C	167	LEU	2.4
1	G	133	ILE	2.4
1	G	346	PHE	2.4
1	E	181	ALA	2.4
1	F	60	LEU	2.4
1	D	322	ASN	2.3
1	G	71	ALA	2.3
1	C	296	PHE	2.3
1	E	296	PHE	2.3
1	H	104	ILE	2.3
1	A	68	LEU	2.3
1	E	73	ALA	2.3
1	G	76	ALA	2.3
1	C	44	ILE	2.3
1	E	274	VAL	2.3
1	A	106	ASP	2.3
1	G	177	ASP	2.3
1	G	322	ASN	2.3
1	F	62	ALA	2.3
1	G	281	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	51	ALA	2.3
1	D	1	MET	2.2
1	H	323	GLY	2.2
1	F	346	PHE	2.2
1	B	177	ASP	2.2
1	C	139	VAL	2.2
1	F	106	ASP	2.2
1	E	77	ILE	2.2
1	E	148	ILE	2.2
1	H	341	VAL	2.2
1	D	105	GLY	2.2
1	C	80	THR	2.2
1	C	131	LEU	2.1
1	D	342	LEU	2.1
1	A	170	ASN	2.1
1	C	49	ASP	2.1
1	E	101	VAL	2.1
1	E	134	ASN	2.1
1	F	53	PHE	2.1
1	C	62	ALA	2.1
1	E	50	SER	2.1
1	C	46	LEU	2.1
1	E	39	MET	2.1
1	F	26	ILE	2.1
1	E	106	ASP	2.1
1	E	168	ASP	2.1
1	H	349	ARG	2.1
1	D	103	ALA	2.1
1	F	174	SER	2.1
1	E	57	LYS	2.1
1	G	163	LEU	2.1
1	G	58	ASN	2.1
1	H	325	ALA	2.1
1	F	319	MET	2.1
1	C	22	GLN	2.1
1	E	46	LEU	2.1
1	E	141	LEU	2.1
1	C	106	ASP	2.1
1	F	170	ASN	2.1
1	G	14	ILE	2.1
1	C	178	VAL	2.1
1	E	94	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	18	GLY	2.1
1	A	91	THR	2.0
1	E	63	GLN	2.0
1	B	20	LEU	2.0
1	C	140	LEU	2.0
1	D	44	ILE	2.0
1	E	130	PRO	2.0
1	E	227	ARG	2.0
1	E	5	GLY	2.0
1	E	68	LEU	2.0
1	E	118	ASP	2.0
1	G	140	LEU	2.0
1	E	14	ILE	2.0
1	F	148	ILE	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	EDO	G	402	4/4	0.60	0.24	53,56,61,63	0
5	SO4	B	407	5/5	0.62	0.14	79,83,87,105	0
5	SO4	F	406	5/5	0.62	0.13	87,101,105,114	0
8	TRS	F	404	8/8	0.63	0.20	59,66,69,76	0
3	EDO	C	405	4/4	0.65	0.17	63,68,69,75	0
2	PEG	A	401	7/7	0.70	0.22	47,50,61,63	0
3	EDO	C	403	4/4	0.74	0.26	55,57,57,62	0
3	EDO	B	403	4/4	0.74	0.21	56,57,58,61	0
3	EDO	E	401	4/4	0.75	0.14	54,63,64,66	0

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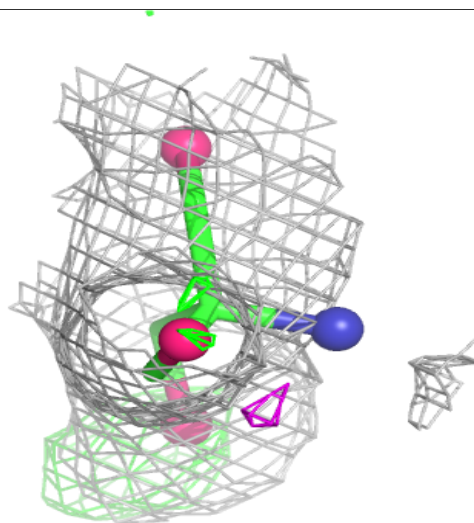
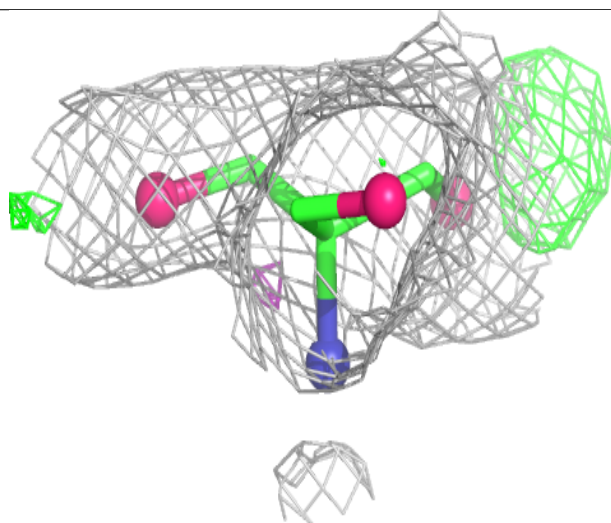
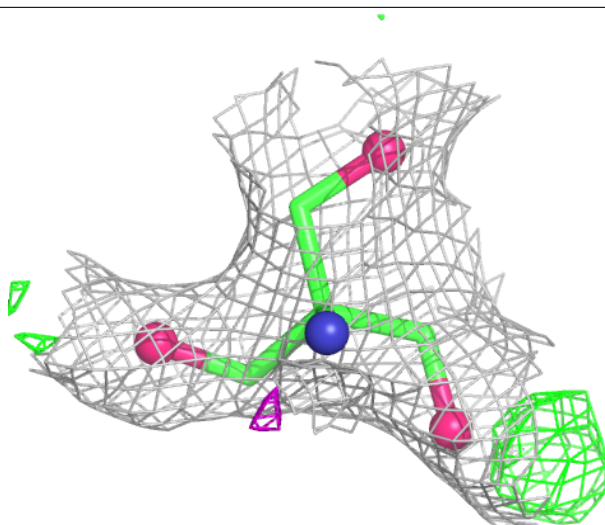
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	F	401	4/4	0.76	0.25	52,52,55,58	0
3	EDO	F	402	4/4	0.76	0.16	52,52,53,56	0
2	PEG	H	401	7/7	0.77	0.17	48,56,72,73	0
3	EDO	B	406	4/4	0.78	0.17	52,54,58,58	0
3	EDO	B	405	4/4	0.78	0.19	45,48,52,53	0
3	EDO	B	404	4/4	0.79	0.14	63,64,69,72	0
3	EDO	H	404	4/4	0.81	0.16	51,51,53,59	0
4	CL	C	406	1/1	0.81	0.16	76,76,76,76	0
2	PEG	C	401	7/7	0.81	0.15	39,51,61,66	0
3	EDO	G	404	4/4	0.81	0.19	50,55,57,58	0
3	EDO	G	405	4/4	0.81	0.15	49,51,51,56	0
3	EDO	A	402	4/4	0.82	0.18	55,58,61,64	0
4	CL	F	405	1/1	0.83	0.20	72,72,72,72	0
3	EDO	C	404	4/4	0.83	0.15	47,48,52,53	0
2	PEG	B	401	7/7	0.84	0.18	53,58,65,66	0
3	EDO	H	402	4/4	0.84	0.16	43,45,49,49	0
4	CL	H	407	1/1	0.85	0.12	73,73,73,73	0
4	CL	H	406	1/1	0.85	0.11	69,69,69,69	0
3	EDO	D	401	4/4	0.86	0.14	57,58,58,60	0
3	EDO	F	403	4/4	0.86	0.15	42,46,48,55	0
3	EDO	C	402	4/4	0.86	0.14	47,47,48,52	0
3	EDO	G	403	4/4	0.86	0.15	50,51,54,57	0
3	EDO	A	404	4/4	0.86	0.14	44,50,52,53	0
3	EDO	H	405	4/4	0.87	0.13	48,49,49,49	0
3	EDO	A	403	4/4	0.87	0.12	48,48,51,52	0
6	GOL	D	404	6/6	0.88	0.20	46,47,52,56	0
3	EDO	G	401	4/4	0.88	0.15	53,58,59,61	0
3	EDO	B	402	4/4	0.90	0.10	44,50,50,51	0
4	CL	D	403	1/1	0.90	0.12	61,61,61,61	0
4	CL	A	405	1/1	0.90	0.14	62,62,62,62	0
3	EDO	H	403	4/4	0.91	0.12	43,44,50,53	0
4	CL	G	406	1/1	0.91	0.12	71,71,71,71	0
3	EDO	D	402	4/4	0.91	0.11	43,46,49,54	0
4	CL	E	402	1/1	0.91	0.13	71,71,71,71	0
7	CA	G	407	1/1	0.94	0.08	64,64,64,64	0
7	CA	D	405	1/1	0.95	0.07	59,59,59,59	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.

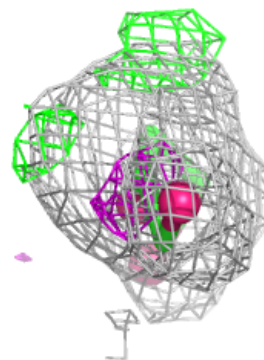
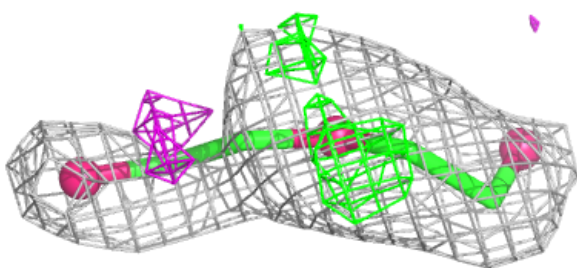
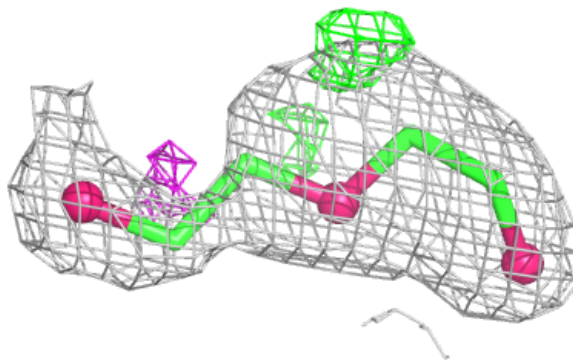
Electron density around TRS F 404:

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

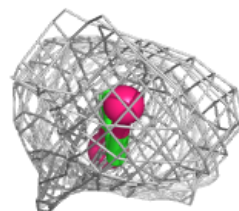
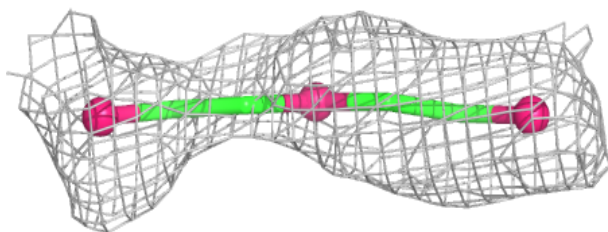
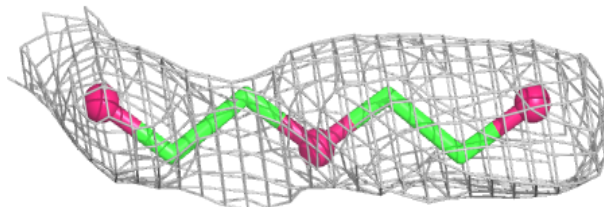


Electron density around PEG A 401:

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

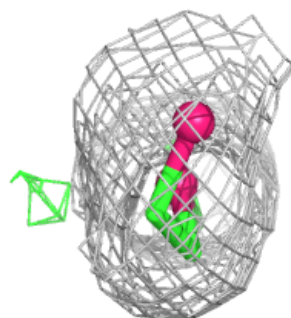
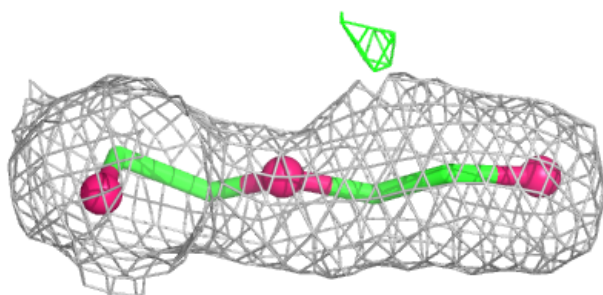
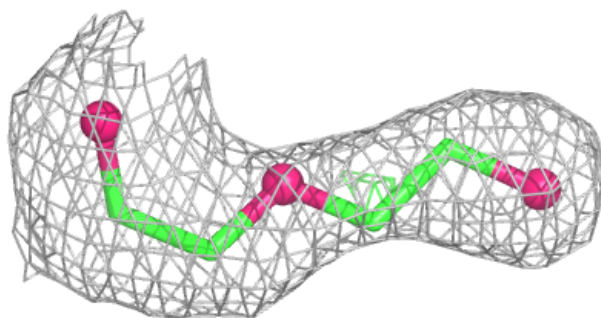
**Electron density around PEG H 401:**

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

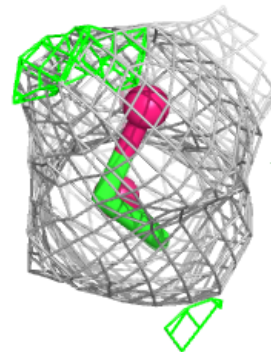
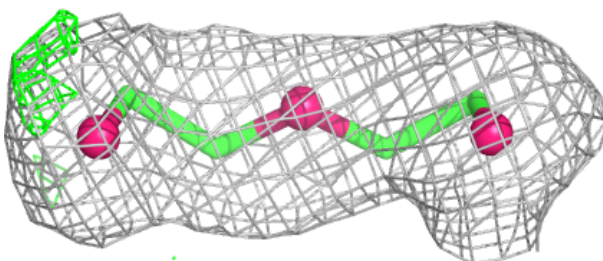
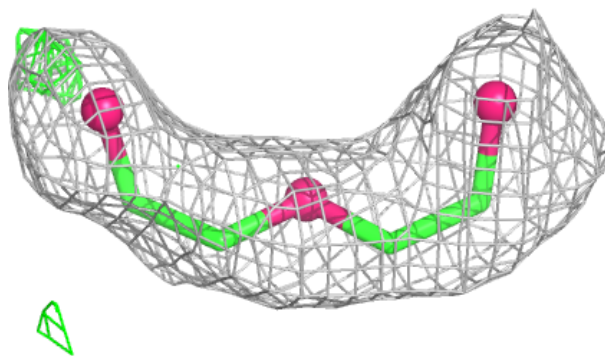


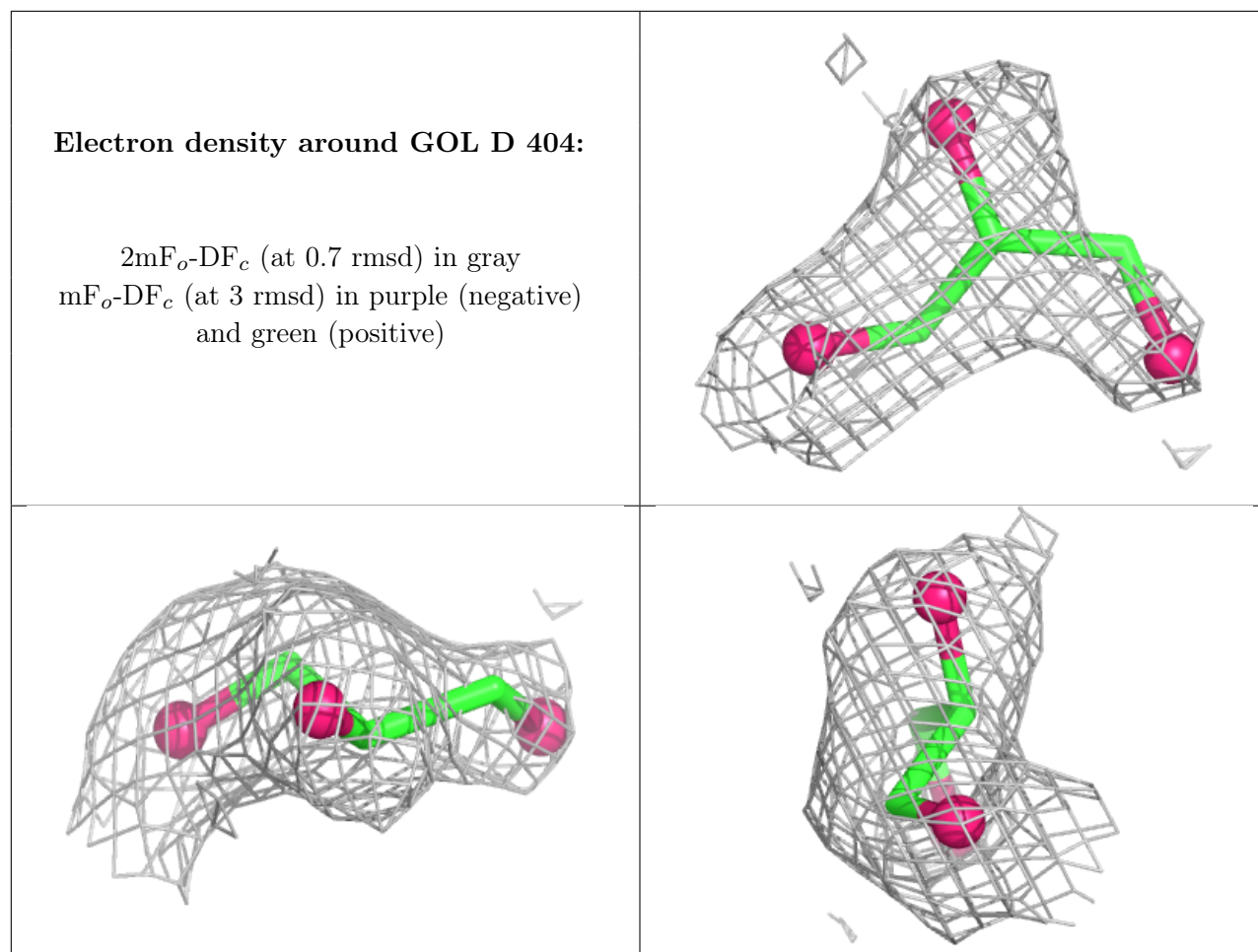
Electron density around PEG C 401:

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

**Electron density around PEG B 401:**

$2mF_o-DF_c$ (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.