



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 9, 2020 – 04:11 PM BST

PDB ID : 6W1K
Title : Crystal structure of the hydroxyglutarate synthase in complex with 2-oxoadipate from *Oryza sativa*
Authors : Pereira, J.H.; Thompson, M.G.; Blake-Hedges, J.M.; Keasling, J.D.; Adams, P.D.
Deposited on : 2020-03-04
Resolution : 1.85 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

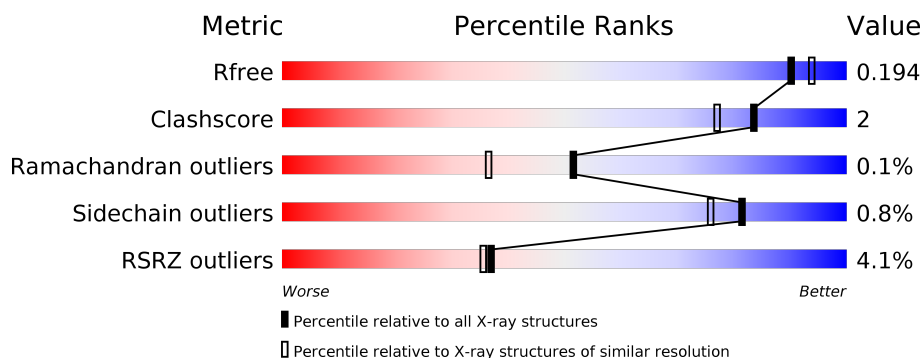
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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

Mol	Chain	Length	Quality of chain
1	A	322	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>7%</div> <div>•</div> </div> </div>
1	B	322	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>7%</div> <div>5%</div> </div> </div>
1	C	322	<div> <div>7%</div> <div> <div></div> <div>90%</div> <div>•</div> <div>5%</div> </div> </div>
1	D	322	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20597 atoms, of which 9546 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 Hydroxyglutarate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	312	Total	C	H	N	O	S	0	0	0
			4865	1573	2395	418	477	2			
1	B	306	Total	C	H	N	O	S	0	1	0
			4818	1556	2373	415	472	2			
1	C	305	Total	C	H	N	O	S	0	1	0
			4820	1556	2376	417	469	2			
1	D	305	Total	C	H	N	O	S	0	1	0
			4811	1554	2370	415	470	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	318	HIS	-	expression tag	UNP Q8H916
A	319	HIS	-	expression tag	UNP Q8H916
A	320	HIS	-	expression tag	UNP Q8H916
A	321	HIS	-	expression tag	UNP Q8H916
A	322	HIS	-	expression tag	UNP Q8H916
A	323	HIS	-	expression tag	UNP Q8H916
B	318	HIS	-	expression tag	UNP Q8H916
B	319	HIS	-	expression tag	UNP Q8H916
B	320	HIS	-	expression tag	UNP Q8H916
B	321	HIS	-	expression tag	UNP Q8H916
B	322	HIS	-	expression tag	UNP Q8H916
B	323	HIS	-	expression tag	UNP Q8H916
C	318	HIS	-	expression tag	UNP Q8H916
C	319	HIS	-	expression tag	UNP Q8H916
C	320	HIS	-	expression tag	UNP Q8H916
C	321	HIS	-	expression tag	UNP Q8H916
C	322	HIS	-	expression tag	UNP Q8H916
C	323	HIS	-	expression tag	UNP Q8H916
D	318	HIS	-	expression tag	UNP Q8H916
D	319	HIS	-	expression tag	UNP Q8H916
D	320	HIS	-	expression tag	UNP Q8H916

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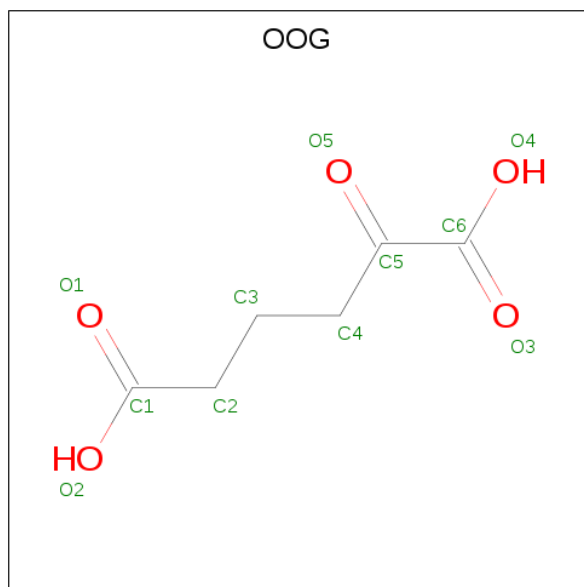
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Chain	Residue	Modelled	Actual	Comment	Reference
D	321	HIS	-	expression tag	UNP Q8H916
D	322	HIS	-	expression tag	UNP Q8H916
D	323	HIS	-	expression tag	UNP Q8H916

- Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ni 1 1	0	0
2	A	1	Total Ni 1 1	0	0
2	D	1	Total Ni 1 1	0	0
2	C	1	Total Ni 1 1	0	0

- Molecule 3 is 2-OXOADIPIC ACID (three-letter code: OOG) (formula: C₆H₈O₅) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 19 6 8 5	0	0
3	B	1	Total C H O 19 6 8 5	0	0
3	C	1	Total C H O 19 6 8 5	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	D	1	Total	C	H	O	0	0
			19	6	8	5		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	286	Total	O	0	0
			286	286		
5	B	283	Total	O	0	0
			283	283		
5	C	242	Total	O	0	0
			242	242		
5	D	312	Total	O	0	0
			312	312		

- Molecule 1: Hydroxyglutarate synthase



HIS	ALA
HIS	ALA
HIS	ALA
HIS	SFR
HIS	ALA
	PHQ
	PHQ
	ALA
	P11
	A12
	R37
	F79
	L128
	L133
	K143
	T147
	T156
	L157
	L177
	S178
	R179
	L188
	A193
	T201
	I211
	L239
	A255
	E268
	F269
	D280
	D284
	R292
	N309
	T313
	R314
	R315
	S316
	ALA
	HIS

P276	ALA	ALA	ALA	ALA	SER	ALA	PRO	ALA	P11	P12	P13	S21	T25	F79	R99	T112	L128	L133	T156	Y171	V176	L177	S178	R179	L188	N195	S210	K216	E228	K233	T253	D256	G257	L258	T259	E260	S265	L273
	ALA	ALA	ALA	ALA	SER	ALA	PRO	ALA	P11	P12	P13	S21	T25	F79	R99	T112	L128	L133	T156	Y171	V176	L177	S178	R179	L188	N195	S210	K216	E228	K233	T253	D256	G257	L258	T259	E260	S265	L273

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.30 Å 98.20 Å 203.80 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.79 – 1.85 42.40 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.0 (39.79-1.85) 99.0 (42.40-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.86 Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.165 , 0.192 0.166 , 0.194	Depositor DCC
R_{free} test set	7756 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	20597	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: NI, OOG, SO4

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	0.55	2/2532 (0.1%)	0.66	1/3436 (0.0%)
1	B	0.63	3/2508 (0.1%)	0.74	6/3398 (0.2%)
1	C	0.54	0/2507	0.64	0/3396
1	D	0.56	0/2504	0.70	4/3393 (0.1%)
All	All	0.57	5/10051 (0.0%)	0.69	11/13623 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	179	ARG	CZ-NH2	-10.42	1.19	1.33
1	B	179	ARG	CZ-NH1	9.00	1.44	1.33
1	B	179	ARG	NE-CZ	-8.09	1.22	1.33
1	A	235	SER	C-N	7.24	1.48	1.34
1	A	235	SER	CA-CB	-5.62	1.44	1.52

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	ARG	NE-CZ-NH2	9.77	125.19	120.30
1	B	179	ARG	NE-CZ-NH1	-7.18	116.71	120.30
1	B	179	ARG	CD-NE-CZ	7.18	133.65	123.60
1	D	99	ARG	NE-CZ-NH1	6.90	123.75	120.30
1	D	99	ARG	NE-CZ-NH2	-6.88	116.86	120.30

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	2470	2395	2395	11	0
1	B	2445	2373	2375	12	0
1	C	2444	2376	2378	8	0
1	D	2441	2370	2371	16	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	11	8	7	0	0
3	B	11	8	7	0	0
3	C	11	8	7	0	0
3	D	11	8	7	0	0
4	A	10	0	0	0	0
4	B	25	0	0	0	0
4	C	25	0	0	1	0
4	D	20	0	0	0	0
5	A	286	0	0	0	0
5	B	283	0	0	1	0
5	C	242	0	0	2	0
5	D	312	0	0	3	0
All	All	11051	9546	9547	46	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 2.

The worst 5 of 46 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:D:176:VAL:HA	5:D:505:HOH:O	1.27	1.30
1:C:150:LYS:NZ	5:C:501:HOH:O	2.03	0.90
1:B:201:THR:HG21	1:B:211:ILE:HD12	1.67	0.77
1:C:309:ASN:O	1:C:313:THR:HG23	1.86	0.76
1:B:316:SER:OG	1:B:316:SER:O	2.05	0.72

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	310/322 (96%)	302 (97%)	8 (3%)	0	100	100
1	B	305/322 (95%)	298 (98%)	6 (2%)	1 (0%)	41	26
1	C	304/322 (94%)	298 (98%)	6 (2%)	0	100	100
1	D	304/322 (94%)	301 (99%)	3 (1%)	0	100	100
All	All	1223/1288 (95%)	1199 (98%)	23 (2%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	315	ARG

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	268/276 (97%)	267 (100%)	1 (0%)	91	89
1	B	268/276 (97%)	267 (100%)	1 (0%)	91	89
1	C	267/276 (97%)	265 (99%)	2 (1%)	84	79
1	D	267/276 (97%)	262 (98%)	5 (2%)	57	43
All	All	1070/1104 (97%)	1061 (99%)	9 (1%)	81	76

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	179	ARG
1	D	292	ARG
1	D	216	LYS
1	C	146	ARG
1	D	210	SER

Some 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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 4 are monoatomic - leaving 20 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	$\# Z > 2$
4	SO4	D	405	-	4,4,4	0.15	0	6,6,6	0.13	0
4	SO4	B	406	-	4,4,4	0.13	0	6,6,6	0.14	0
4	SO4	C	401	-	4,4,4	0.14	0	6,6,6	0.10	0
4	SO4	A	403	-	4,4,4	0.13	0	6,6,6	0.07	0
4	SO4	B	407	-	4,4,4	0.15	0	6,6,6	0.07	0
3	OOG	D	402	2	4,10,10	1.51	1 (25%)	4,12,12	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	OOG	A	402	2	4,10,10	1.48	1 (25%)	4,12,12	0.87	0
3	OOG	B	402	2	4,10,10	1.38	1 (25%)	4,12,12	0.97	0
4	SO4	C	407	-	4,4,4	0.13	0	6,6,6	0.16	0
4	SO4	D	403	-	4,4,4	0.17	0	6,6,6	0.19	0
4	SO4	C	406	-	4,4,4	0.16	0	6,6,6	0.09	0
4	SO4	D	404	-	4,4,4	0.18	0	6,6,6	0.37	0
4	SO4	B	404	-	4,4,4	0.12	0	6,6,6	0.13	0
4	SO4	B	403	-	4,4,4	0.15	0	6,6,6	0.09	0
3	OOG	C	404	2	4,10,10	1.41	1 (25%)	4,12,12	1.03	0
4	SO4	C	405	-	4,4,4	0.13	0	6,6,6	0.10	0
4	SO4	C	402	-	4,4,4	0.98	0	6,6,6	0.26	0
4	SO4	B	405	-	4,4,4	0.17	0	6,6,6	0.07	0
4	SO4	D	406	-	4,4,4	0.14	0	6,6,6	0.08	0
4	SO4	A	404	-	4,4,4	0.14	0	6,6,6	0.10	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	OOG	D	402	2	-	2/4/10/10	-
3	OOG	A	402	2	-	2/4/10/10	-
3	OOG	B	402	2	-	2/4/10/10	-
3	OOG	C	404	2	-	2/4/10/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	402	OOG	O5-C5	-2.54	1.18	1.22
3	C	404	OOG	O5-C5	-2.46	1.18	1.22
3	A	402	OOG	O5-C5	-2.45	1.18	1.22
3	B	402	OOG	O5-C5	-2.38	1.18	1.22

There are no bond angle outliers.

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	404	OOG	C3-C4-C5-C6

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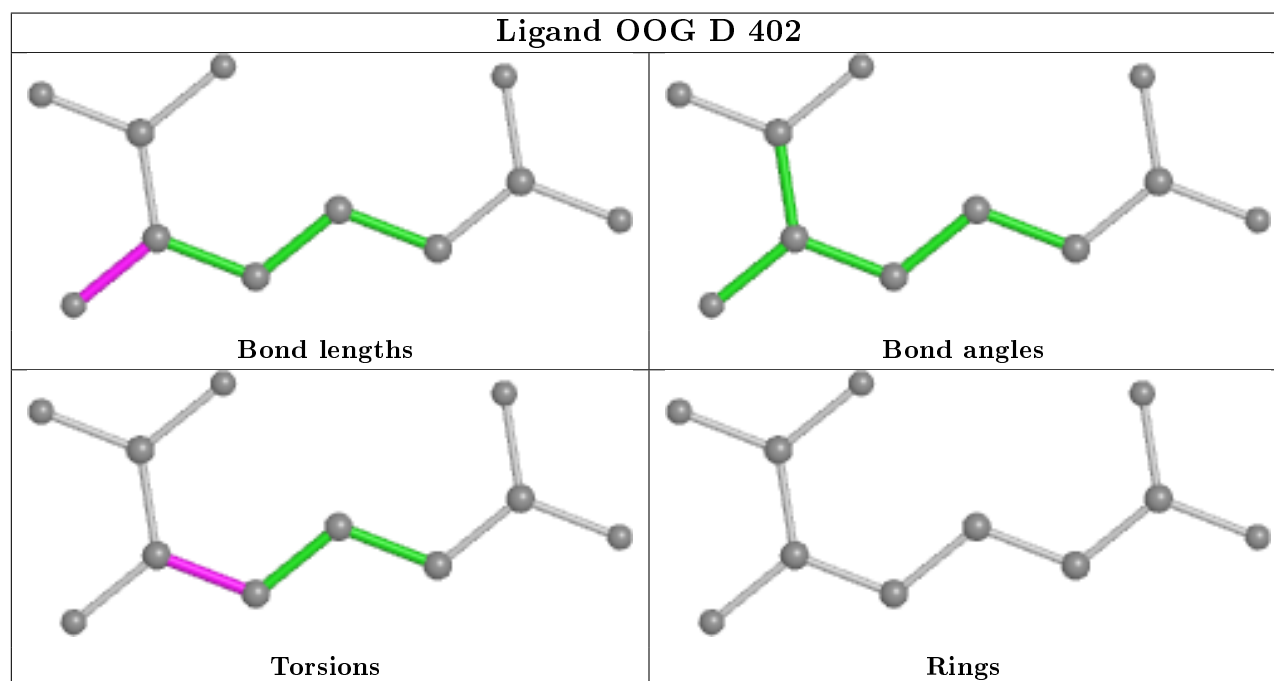
Mol	Chain	Res	Type	Atoms
3	B	402	OOG	C3-C4-C5-C6
3	C	404	OOG	C3-C4-C5-O5
3	B	402	OOG	C3-C4-C5-O5
3	A	402	OOG	C3-C4-C5-C6

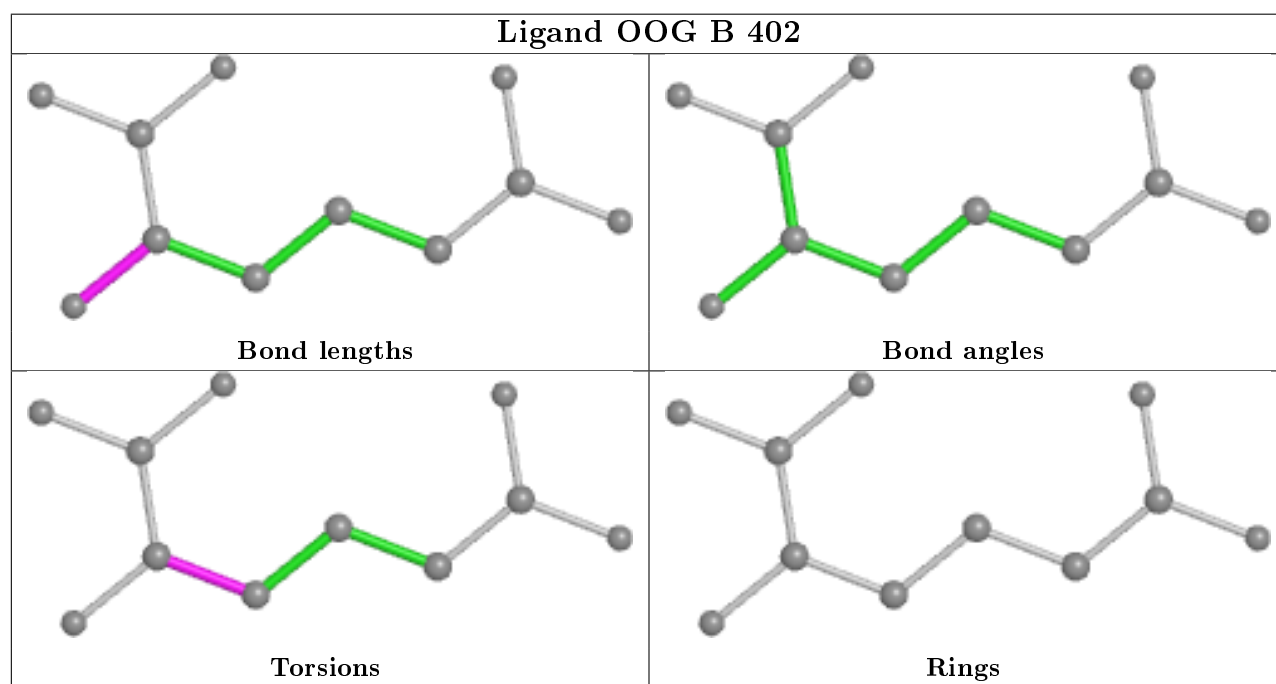
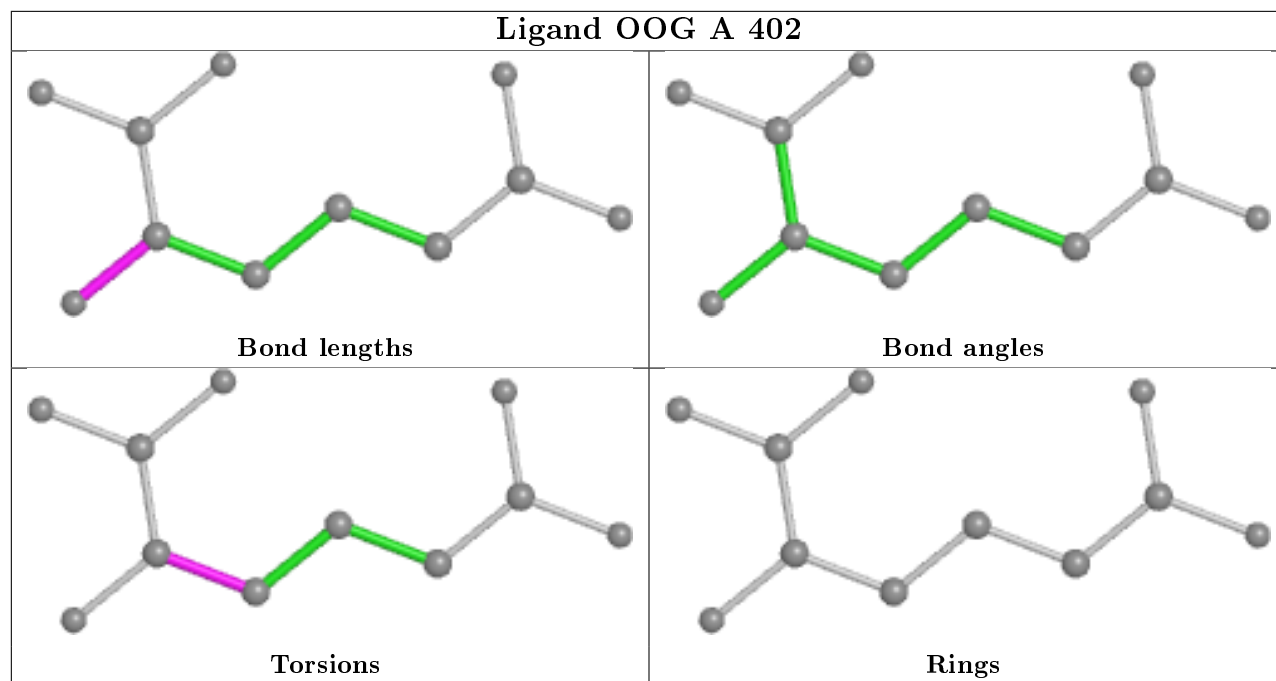
There are no ring outliers.

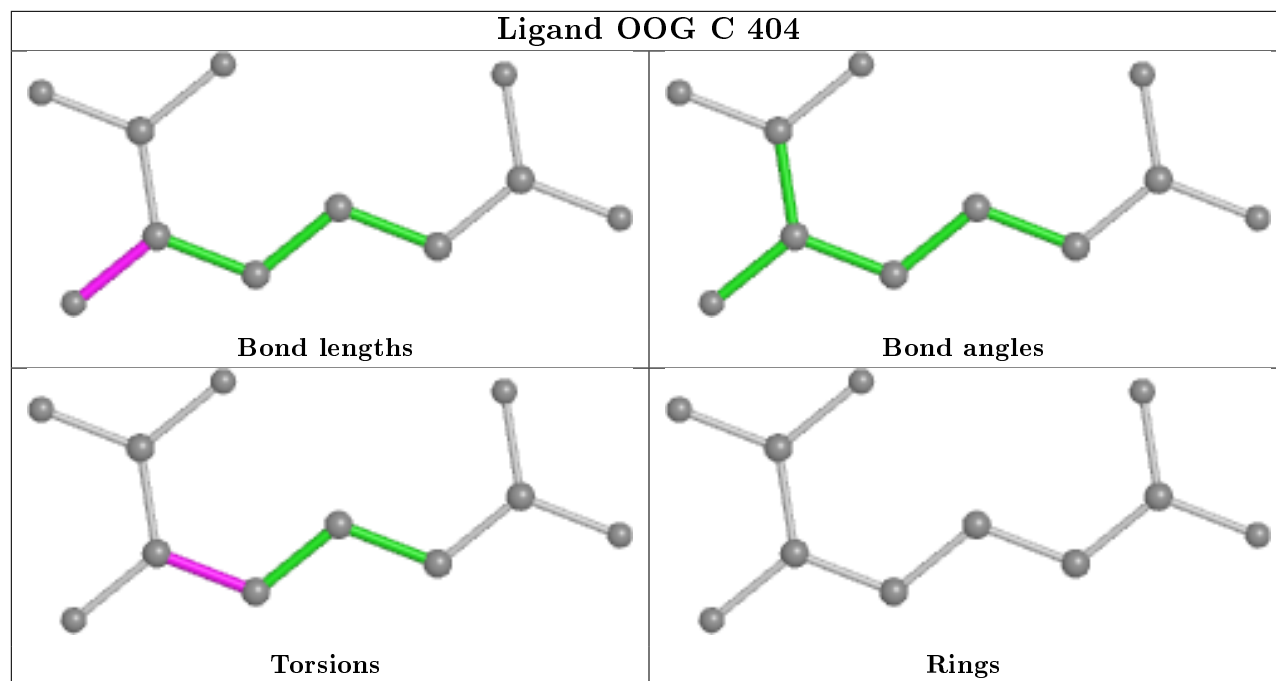
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	405	SO4	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 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 [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	312/322 (96%)	0.32	13 (4%) 36 34	21, 35, 68, 109	0
1	B	306/322 (95%)	0.05	9 (2%) 51 50	22, 32, 60, 192	0
1	C	305/322 (94%)	0.35	22 (7%) 15 15	23, 37, 66, 86	0
1	D	305/322 (94%)	0.07	6 (1%) 65 64	20, 30, 59, 99	0
All	All	1228/1288 (95%)	0.20	50 (4%) 37 35	20, 34, 64, 192	0

The worst 5 of 50 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	282	PRO	8.6
1	A	280	ASP	6.8
1	A	281	LEU	5.6
1	D	12	ALA	5.3
1	C	281	LEU	5.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

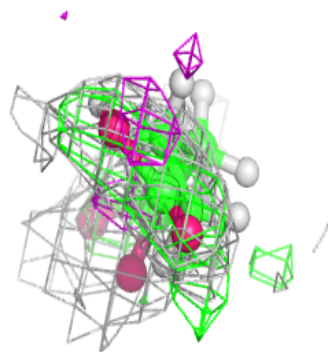
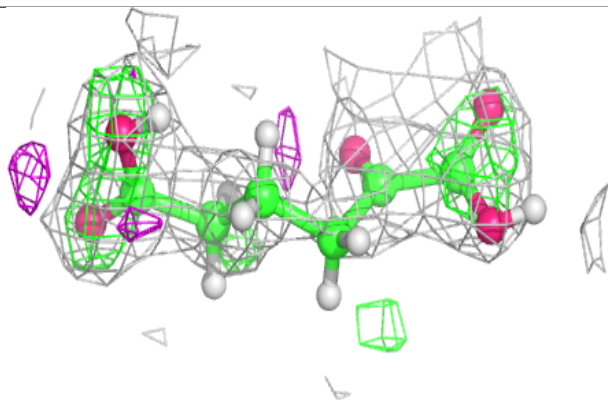
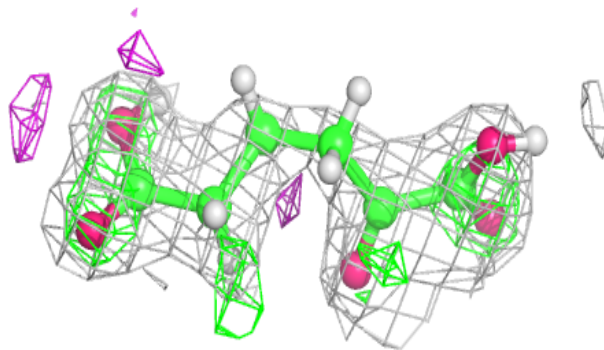
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	B	407	5/5	0.74	0.23	113,113,113,113	0
4	SO4	C	406	5/5	0.76	0.26	98,98,99,99	0
4	SO4	B	404	5/5	0.80	0.22	100,100,101,101	0
4	SO4	A	404	5/5	0.81	0.28	113,113,113,113	0
3	OOG	B	402	11/11	0.82	0.35	96,101,121,121	0
4	SO4	D	406	5/5	0.83	0.29	115,115,116,116	0
4	SO4	B	403	5/5	0.84	0.16	100,100,100,100	0
3	OOG	C	404	11/11	0.87	0.32	83,87,105,105	0
4	SO4	D	405	5/5	0.88	0.16	99,99,99,99	0
4	SO4	C	401	5/5	0.89	0.18	111,111,111,111	0
3	OOG	A	402	11/11	0.89	0.26	76,78,95,95	0
3	OOG	D	402	11/11	0.91	0.13	46,51,60,61	0
4	SO4	D	404	5/5	0.93	0.14	83,83,84,85	0
4	SO4	A	403	5/5	0.93	0.23	104,105,105,105	0
4	SO4	C	402	5/5	0.94	0.17	88,88,89,91	0
4	SO4	C	407	5/5	0.94	0.42	106,106,106,106	0
4	SO4	B	406	5/5	0.94	0.28	111,111,111,111	0
4	SO4	C	405	5/5	0.95	0.09	95,95,96,96	0
4	SO4	B	405	5/5	0.97	0.16	81,81,82,82	0
4	SO4	D	403	5/5	0.97	0.08	66,67,67,68	0
2	NI	B	401	1/1	0.99	0.09	33,33,33,33	1
2	NI	D	401	1/1	1.00	0.08	32,32,32,32	1
2	NI	C	403	1/1	1.00	0.09	36,36,36,36	1
2	NI	A	401	1/1	1.00	0.09	36,36,36,36	1

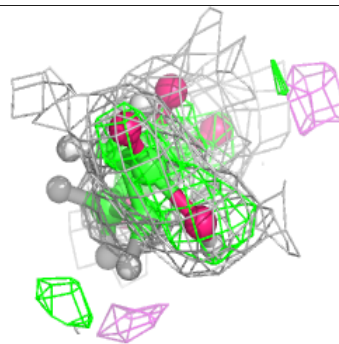
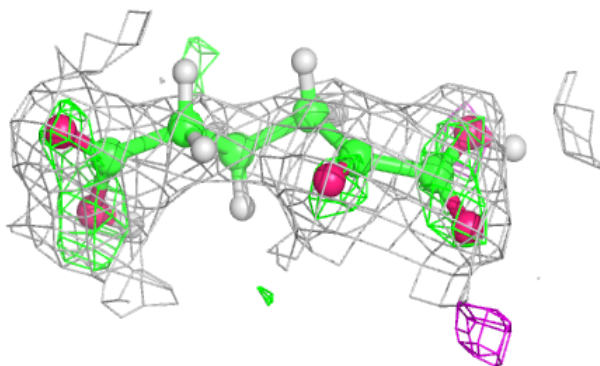
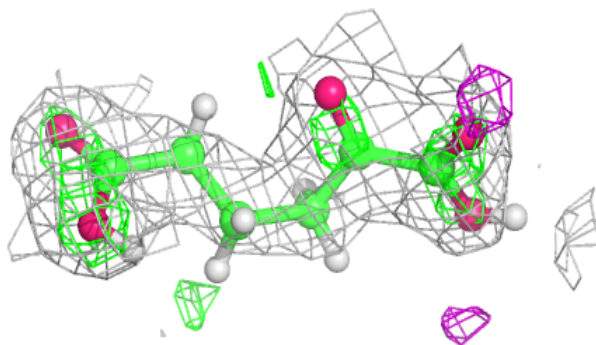
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OOG B 402:

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

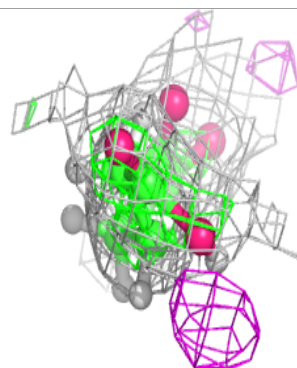
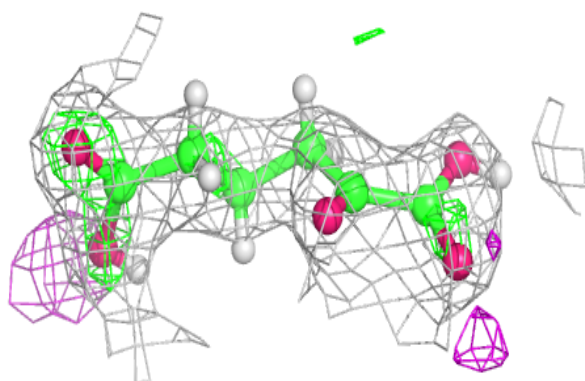
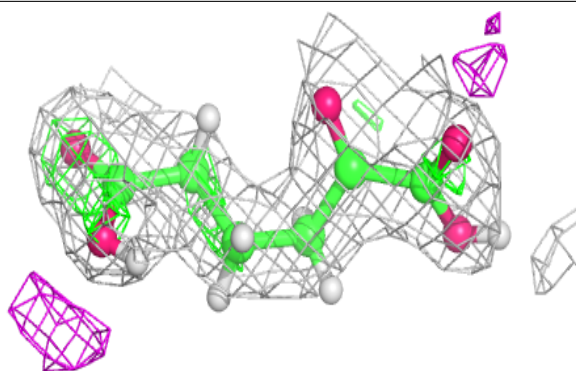
**Electron density around OOG C 404:**

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

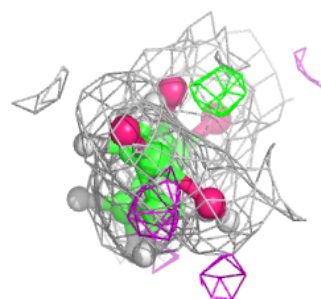
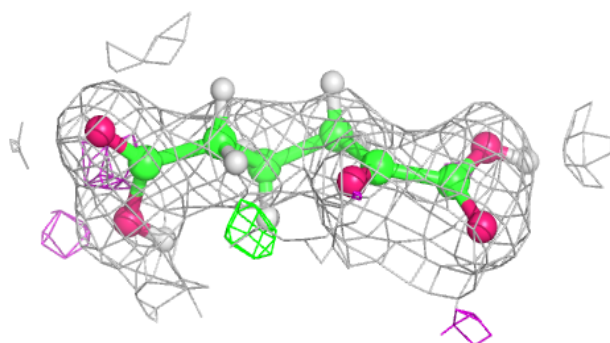
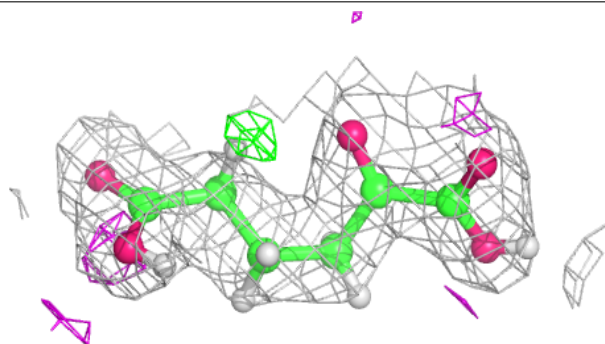


Electron density around OOG A 402:

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

**Electron density around OOG D 402:**

$2mF_o-DF_c$ (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.