



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:26 AM UTC

PDB ID : 5TJR / pdb_00005tjr
Title : X-ray Crystal structure of a methylmalonate semialdehyde dehydrogenase from Pseudomonas sp. AAC
Authors : Peat, T.S.; Newman, J.
Deposited on : 2016-10-05
Resolution : 2.95 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:

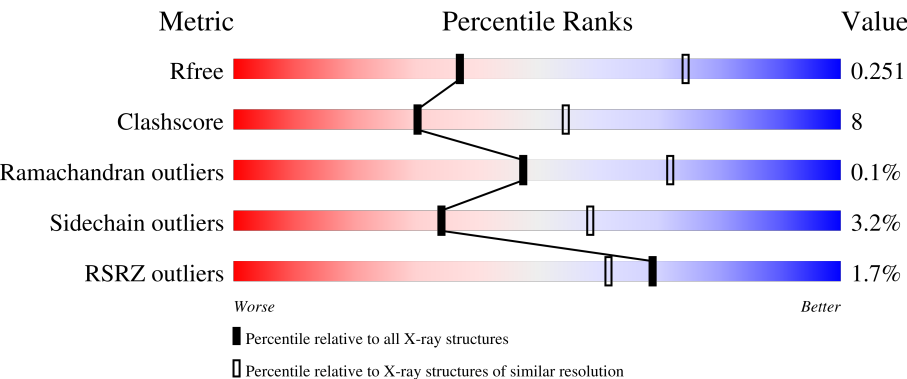
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	531	% 71%14%•13%
1	B	531	% 70%16%•12%
1	C	531	3% 70%16%•13%
1	D	531	% 70%16%•12%
1	E	531	% 71%14%•14%

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Mol	Chain	Length	Quality of chain
1	F	531	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20975 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 Methylmalonate-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	S	0	0	0
			3471	2210	600	646	15			
1	B	465	Total	C	N	O	S	0	1	0
			3525	2249	607	654	15			
1	C	461	Total	C	N	O	S	0	0	0
			3455	2196	601	643	15			
1	D	468	Total	C	N	O	S	0	0	0
			3535	2252	613	655	15			
1	E	455	Total	C	N	O	S	0	0	0
			3408	2170	590	633	15			
1	F	454	Total	C	N	O	S	0	0	0
			3430	2185	594	636	15			

There are 198 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-32	MET	-	initiating methionine	UNP A0A081YAY7
A	-31	ARG	-	expression tag	UNP A0A081YAY7
A	-30	GLY	-	expression tag	UNP A0A081YAY7
A	-29	SER	-	expression tag	UNP A0A081YAY7
A	-28	HIS	-	expression tag	UNP A0A081YAY7
A	-27	HIS	-	expression tag	UNP A0A081YAY7
A	-26	HIS	-	expression tag	UNP A0A081YAY7
A	-25	HIS	-	expression tag	UNP A0A081YAY7
A	-24	HIS	-	expression tag	UNP A0A081YAY7
A	-23	HIS	-	expression tag	UNP A0A081YAY7
A	-22	GLY	-	expression tag	UNP A0A081YAY7
A	-21	MET	-	expression tag	UNP A0A081YAY7
A	-20	ALA	-	expression tag	UNP A0A081YAY7
A	-19	SER	-	expression tag	UNP A0A081YAY7
A	-18	MET	-	expression tag	UNP A0A081YAY7
A	-17	THR	-	expression tag	UNP A0A081YAY7
A	-16	GLY	-	expression tag	UNP A0A081YAY7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	GLY	-	expression tag	UNP A0A081YAY7
A	-14	GLN	-	expression tag	UNP A0A081YAY7
A	-13	GLN	-	expression tag	UNP A0A081YAY7
A	-12	MET	-	expression tag	UNP A0A081YAY7
A	-11	GLY	-	expression tag	UNP A0A081YAY7
A	-10	ARG	-	expression tag	UNP A0A081YAY7
A	-9	ASP	-	expression tag	UNP A0A081YAY7
A	-8	LEU	-	expression tag	UNP A0A081YAY7
A	-7	TYR	-	expression tag	UNP A0A081YAY7
A	-6	ASP	-	expression tag	UNP A0A081YAY7
A	-5	ASP	-	expression tag	UNP A0A081YAY7
A	-4	ASP	-	expression tag	UNP A0A081YAY7
A	-3	ASP	-	expression tag	UNP A0A081YAY7
A	-2	LYS	-	expression tag	UNP A0A081YAY7
A	-1	GLY	-	expression tag	UNP A0A081YAY7
A	0	SER	-	expression tag	UNP A0A081YAY7
B	-32	MET	-	initiating methionine	UNP A0A081YAY7
B	-31	ARG	-	expression tag	UNP A0A081YAY7
B	-30	GLY	-	expression tag	UNP A0A081YAY7
B	-29	SER	-	expression tag	UNP A0A081YAY7
B	-28	HIS	-	expression tag	UNP A0A081YAY7
B	-27	HIS	-	expression tag	UNP A0A081YAY7
B	-26	HIS	-	expression tag	UNP A0A081YAY7
B	-25	HIS	-	expression tag	UNP A0A081YAY7
B	-24	HIS	-	expression tag	UNP A0A081YAY7
B	-23	HIS	-	expression tag	UNP A0A081YAY7
B	-22	GLY	-	expression tag	UNP A0A081YAY7
B	-21	MET	-	expression tag	UNP A0A081YAY7
B	-20	ALA	-	expression tag	UNP A0A081YAY7
B	-19	SER	-	expression tag	UNP A0A081YAY7
B	-18	MET	-	expression tag	UNP A0A081YAY7
B	-17	THR	-	expression tag	UNP A0A081YAY7
B	-16	GLY	-	expression tag	UNP A0A081YAY7
B	-15	GLY	-	expression tag	UNP A0A081YAY7
B	-14	GLN	-	expression tag	UNP A0A081YAY7
B	-13	GLN	-	expression tag	UNP A0A081YAY7
B	-12	MET	-	expression tag	UNP A0A081YAY7
B	-11	GLY	-	expression tag	UNP A0A081YAY7
B	-10	ARG	-	expression tag	UNP A0A081YAY7
B	-9	ASP	-	expression tag	UNP A0A081YAY7
B	-8	LEU	-	expression tag	UNP A0A081YAY7
B	-7	TYR	-	expression tag	UNP A0A081YAY7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	ASP	-	expression tag	UNP A0A081YAY7
B	-5	ASP	-	expression tag	UNP A0A081YAY7
B	-4	ASP	-	expression tag	UNP A0A081YAY7
B	-3	ASP	-	expression tag	UNP A0A081YAY7
B	-2	LYS	-	expression tag	UNP A0A081YAY7
B	-1	GLY	-	expression tag	UNP A0A081YAY7
B	0	SER	-	expression tag	UNP A0A081YAY7
C	-32	MET	-	initiating methionine	UNP A0A081YAY7
C	-31	ARG	-	expression tag	UNP A0A081YAY7
C	-30	GLY	-	expression tag	UNP A0A081YAY7
C	-29	SER	-	expression tag	UNP A0A081YAY7
C	-28	HIS	-	expression tag	UNP A0A081YAY7
C	-27	HIS	-	expression tag	UNP A0A081YAY7
C	-26	HIS	-	expression tag	UNP A0A081YAY7
C	-25	HIS	-	expression tag	UNP A0A081YAY7
C	-24	HIS	-	expression tag	UNP A0A081YAY7
C	-23	HIS	-	expression tag	UNP A0A081YAY7
C	-22	GLY	-	expression tag	UNP A0A081YAY7
C	-21	MET	-	expression tag	UNP A0A081YAY7
C	-20	ALA	-	expression tag	UNP A0A081YAY7
C	-19	SER	-	expression tag	UNP A0A081YAY7
C	-18	MET	-	expression tag	UNP A0A081YAY7
C	-17	THR	-	expression tag	UNP A0A081YAY7
C	-16	GLY	-	expression tag	UNP A0A081YAY7
C	-15	GLY	-	expression tag	UNP A0A081YAY7
C	-14	GLN	-	expression tag	UNP A0A081YAY7
C	-13	GLN	-	expression tag	UNP A0A081YAY7
C	-12	MET	-	expression tag	UNP A0A081YAY7
C	-11	GLY	-	expression tag	UNP A0A081YAY7
C	-10	ARG	-	expression tag	UNP A0A081YAY7
C	-9	ASP	-	expression tag	UNP A0A081YAY7
C	-8	LEU	-	expression tag	UNP A0A081YAY7
C	-7	TYR	-	expression tag	UNP A0A081YAY7
C	-6	ASP	-	expression tag	UNP A0A081YAY7
C	-5	ASP	-	expression tag	UNP A0A081YAY7
C	-4	ASP	-	expression tag	UNP A0A081YAY7
C	-3	ASP	-	expression tag	UNP A0A081YAY7
C	-2	LYS	-	expression tag	UNP A0A081YAY7
C	-1	GLY	-	expression tag	UNP A0A081YAY7
C	0	SER	-	expression tag	UNP A0A081YAY7
D	-32	MET	-	initiating methionine	UNP A0A081YAY7
D	-31	ARG	-	expression tag	UNP A0A081YAY7

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-30	GLY	-	expression tag	UNP A0A081YAY7
D	-29	SER	-	expression tag	UNP A0A081YAY7
D	-28	HIS	-	expression tag	UNP A0A081YAY7
D	-27	HIS	-	expression tag	UNP A0A081YAY7
D	-26	HIS	-	expression tag	UNP A0A081YAY7
D	-25	HIS	-	expression tag	UNP A0A081YAY7
D	-24	HIS	-	expression tag	UNP A0A081YAY7
D	-23	HIS	-	expression tag	UNP A0A081YAY7
D	-22	GLY	-	expression tag	UNP A0A081YAY7
D	-21	MET	-	expression tag	UNP A0A081YAY7
D	-20	ALA	-	expression tag	UNP A0A081YAY7
D	-19	SER	-	expression tag	UNP A0A081YAY7
D	-18	MET	-	expression tag	UNP A0A081YAY7
D	-17	THR	-	expression tag	UNP A0A081YAY7
D	-16	GLY	-	expression tag	UNP A0A081YAY7
D	-15	GLY	-	expression tag	UNP A0A081YAY7
D	-14	GLN	-	expression tag	UNP A0A081YAY7
D	-13	GLN	-	expression tag	UNP A0A081YAY7
D	-12	MET	-	expression tag	UNP A0A081YAY7
D	-11	GLY	-	expression tag	UNP A0A081YAY7
D	-10	ARG	-	expression tag	UNP A0A081YAY7
D	-9	ASP	-	expression tag	UNP A0A081YAY7
D	-8	LEU	-	expression tag	UNP A0A081YAY7
D	-7	TYR	-	expression tag	UNP A0A081YAY7
D	-6	ASP	-	expression tag	UNP A0A081YAY7
D	-5	ASP	-	expression tag	UNP A0A081YAY7
D	-4	ASP	-	expression tag	UNP A0A081YAY7
D	-3	ASP	-	expression tag	UNP A0A081YAY7
D	-2	LYS	-	expression tag	UNP A0A081YAY7
D	-1	GLY	-	expression tag	UNP A0A081YAY7
D	0	SER	-	expression tag	UNP A0A081YAY7
E	-32	MET	-	initiating methionine	UNP A0A081YAY7
E	-31	ARG	-	expression tag	UNP A0A081YAY7
E	-30	GLY	-	expression tag	UNP A0A081YAY7
E	-29	SER	-	expression tag	UNP A0A081YAY7
E	-28	HIS	-	expression tag	UNP A0A081YAY7
E	-27	HIS	-	expression tag	UNP A0A081YAY7
E	-26	HIS	-	expression tag	UNP A0A081YAY7
E	-25	HIS	-	expression tag	UNP A0A081YAY7
E	-24	HIS	-	expression tag	UNP A0A081YAY7
E	-23	HIS	-	expression tag	UNP A0A081YAY7
E	-22	GLY	-	expression tag	UNP A0A081YAY7

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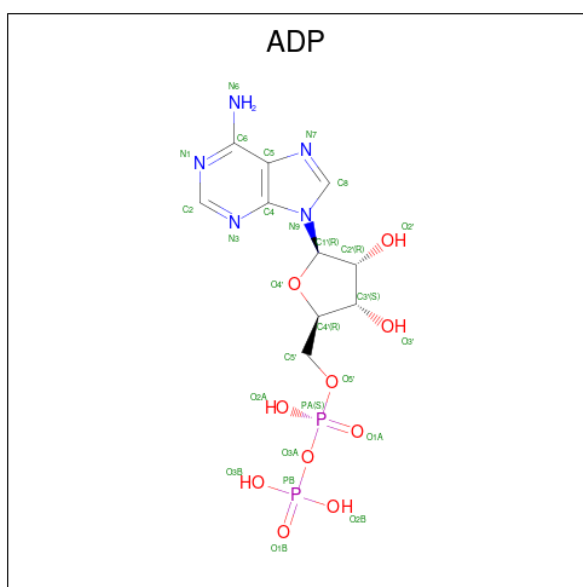
Chain	Residue	Modelled	Actual	Comment	Reference
E	-21	MET	-	expression tag	UNP A0A081YAY7
E	-20	ALA	-	expression tag	UNP A0A081YAY7
E	-19	SER	-	expression tag	UNP A0A081YAY7
E	-18	MET	-	expression tag	UNP A0A081YAY7
E	-17	THR	-	expression tag	UNP A0A081YAY7
E	-16	GLY	-	expression tag	UNP A0A081YAY7
E	-15	GLY	-	expression tag	UNP A0A081YAY7
E	-14	GLN	-	expression tag	UNP A0A081YAY7
E	-13	GLN	-	expression tag	UNP A0A081YAY7
E	-12	MET	-	expression tag	UNP A0A081YAY7
E	-11	GLY	-	expression tag	UNP A0A081YAY7
E	-10	ARG	-	expression tag	UNP A0A081YAY7
E	-9	ASP	-	expression tag	UNP A0A081YAY7
E	-8	LEU	-	expression tag	UNP A0A081YAY7
E	-7	TYR	-	expression tag	UNP A0A081YAY7
E	-6	ASP	-	expression tag	UNP A0A081YAY7
E	-5	ASP	-	expression tag	UNP A0A081YAY7
E	-4	ASP	-	expression tag	UNP A0A081YAY7
E	-3	ASP	-	expression tag	UNP A0A081YAY7
E	-2	LYS	-	expression tag	UNP A0A081YAY7
E	-1	GLY	-	expression tag	UNP A0A081YAY7
E	0	SER	-	expression tag	UNP A0A081YAY7
F	-32	MET	-	initiating methionine	UNP A0A081YAY7
F	-31	ARG	-	expression tag	UNP A0A081YAY7
F	-30	GLY	-	expression tag	UNP A0A081YAY7
F	-29	SER	-	expression tag	UNP A0A081YAY7
F	-28	HIS	-	expression tag	UNP A0A081YAY7
F	-27	HIS	-	expression tag	UNP A0A081YAY7
F	-26	HIS	-	expression tag	UNP A0A081YAY7
F	-25	HIS	-	expression tag	UNP A0A081YAY7
F	-24	HIS	-	expression tag	UNP A0A081YAY7
F	-23	HIS	-	expression tag	UNP A0A081YAY7
F	-22	GLY	-	expression tag	UNP A0A081YAY7
F	-21	MET	-	expression tag	UNP A0A081YAY7
F	-20	ALA	-	expression tag	UNP A0A081YAY7
F	-19	SER	-	expression tag	UNP A0A081YAY7
F	-18	MET	-	expression tag	UNP A0A081YAY7
F	-17	THR	-	expression tag	UNP A0A081YAY7
F	-16	GLY	-	expression tag	UNP A0A081YAY7
F	-15	GLY	-	expression tag	UNP A0A081YAY7
F	-14	GLN	-	expression tag	UNP A0A081YAY7
F	-13	GLN	-	expression tag	UNP A0A081YAY7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-12	MET	-	expression tag	UNP A0A081YAY7
F	-11	GLY	-	expression tag	UNP A0A081YAY7
F	-10	ARG	-	expression tag	UNP A0A081YAY7
F	-9	ASP	-	expression tag	UNP A0A081YAY7
F	-8	LEU	-	expression tag	UNP A0A081YAY7
F	-7	TYR	-	expression tag	UNP A0A081YAY7
F	-6	ASP	-	expression tag	UNP A0A081YAY7
F	-5	ASP	-	expression tag	UNP A0A081YAY7
F	-4	ASP	-	expression tag	UNP A0A081YAY7
F	-3	ASP	-	expression tag	UNP A0A081YAY7
F	-2	LYS	-	expression tag	UNP A0A081YAY7
F	-1	GLY	-	expression tag	UNP A0A081YAY7
F	0	SER	-	expression tag	UNP A0A081YAY7

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

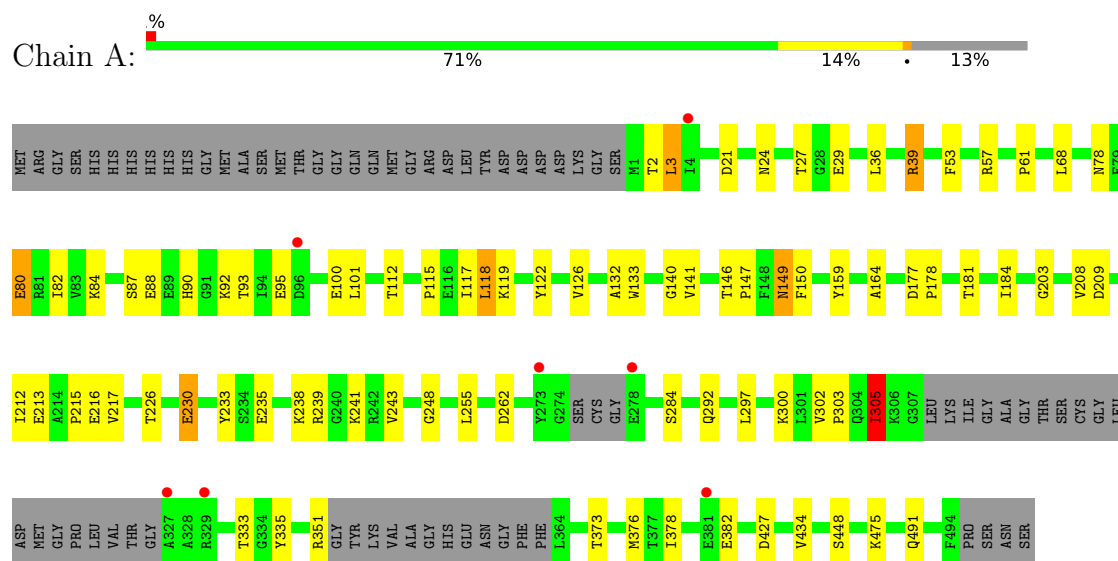
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total 3	O 3	0	0
3	B	5	Total 5	O 5	0	0
3	C	2	Total 2	O 2	0	0
3	D	3	Total 3	O 3	0	0
3	E	1	Total 1	O 1	0	0
3	F	2	Total 2	O 2	0	0

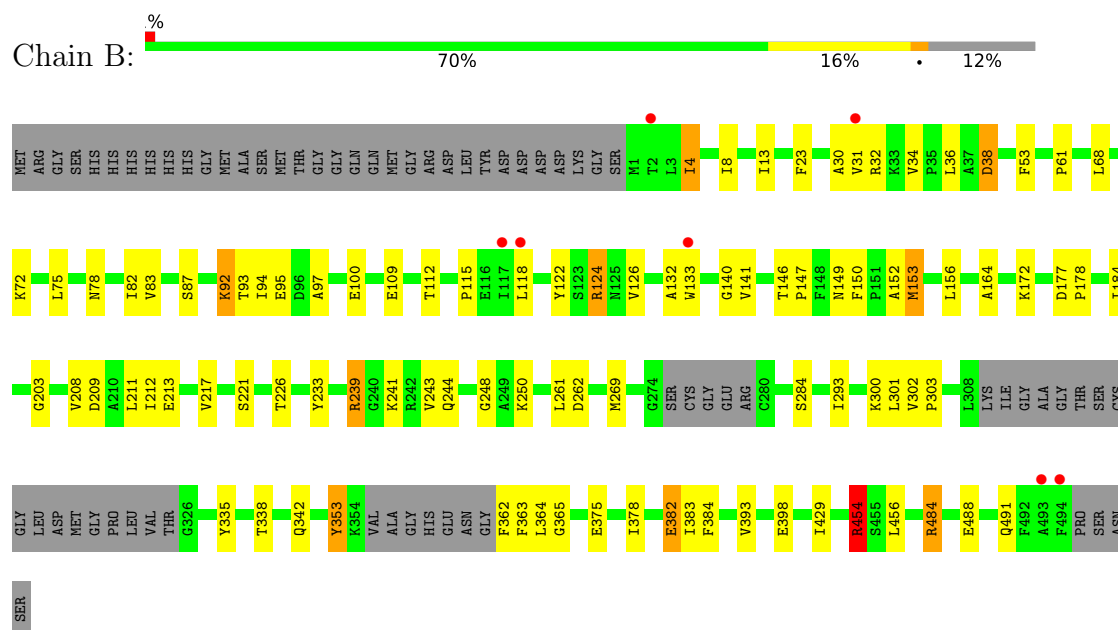
3 Residue-property plots [i](#)

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.

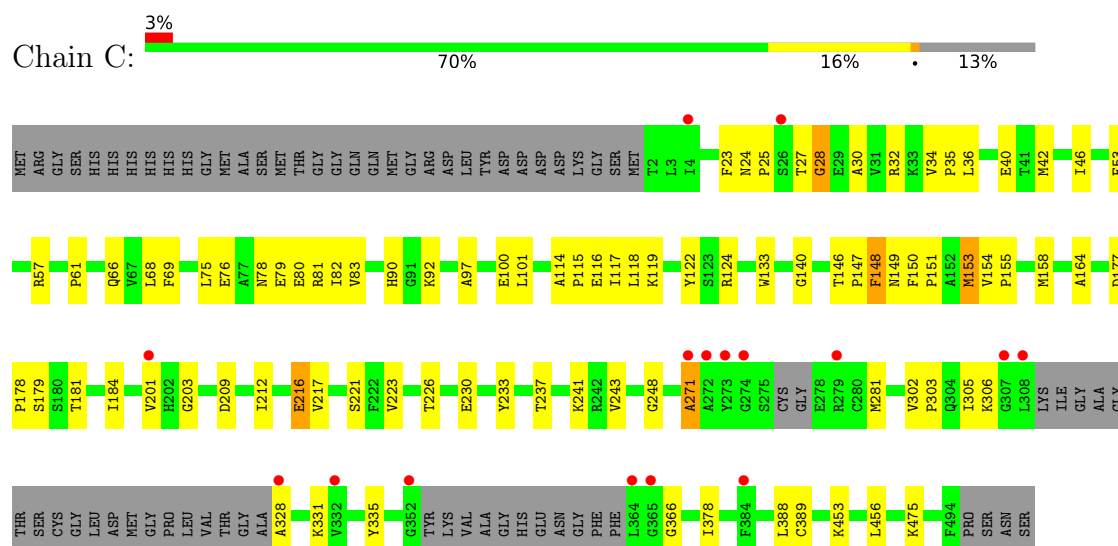
- Molecule 1: Methylmalonate-semialdehyde dehydrogenase



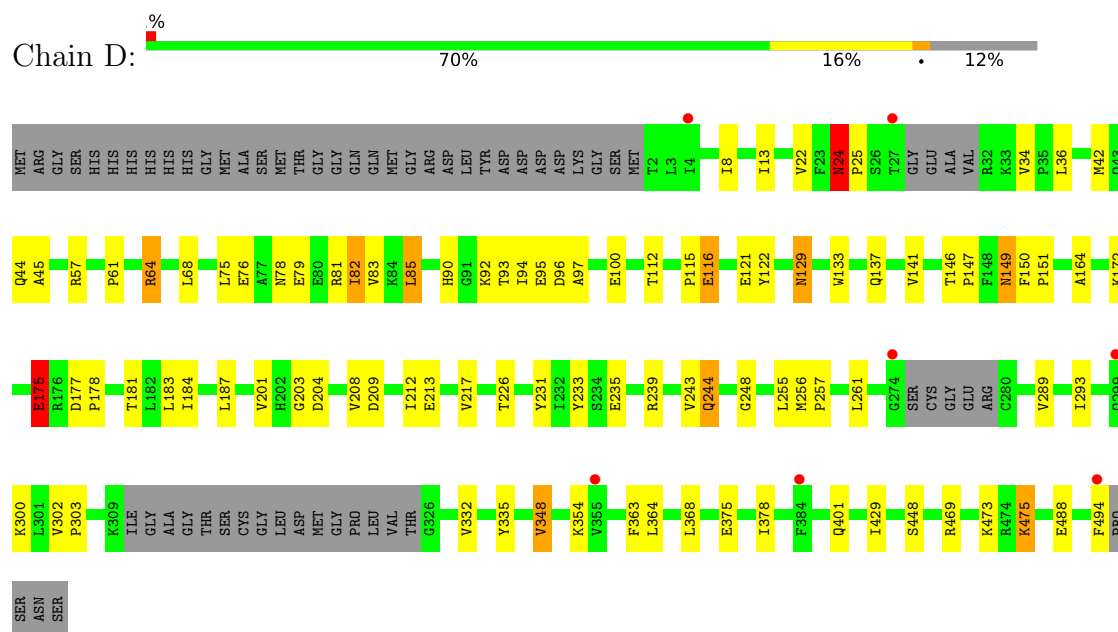
- Molecule 1: Methylmalonate-semialdehyde dehydrogenase



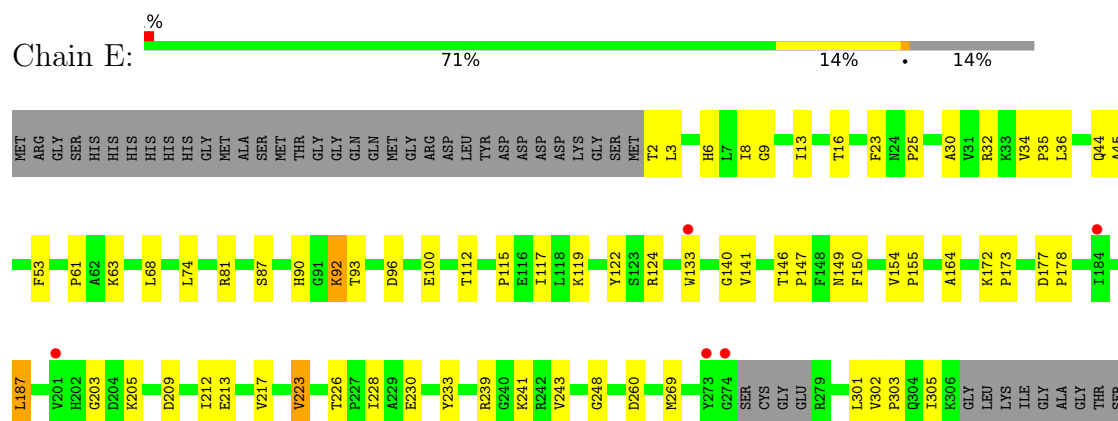
- Molecule 1: Methylmalonate-semialdehyde dehydrogenase

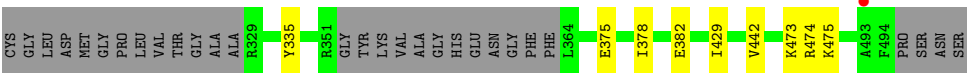


- Molecule 1: Methylmalonate-semialdehyde dehydrogenase

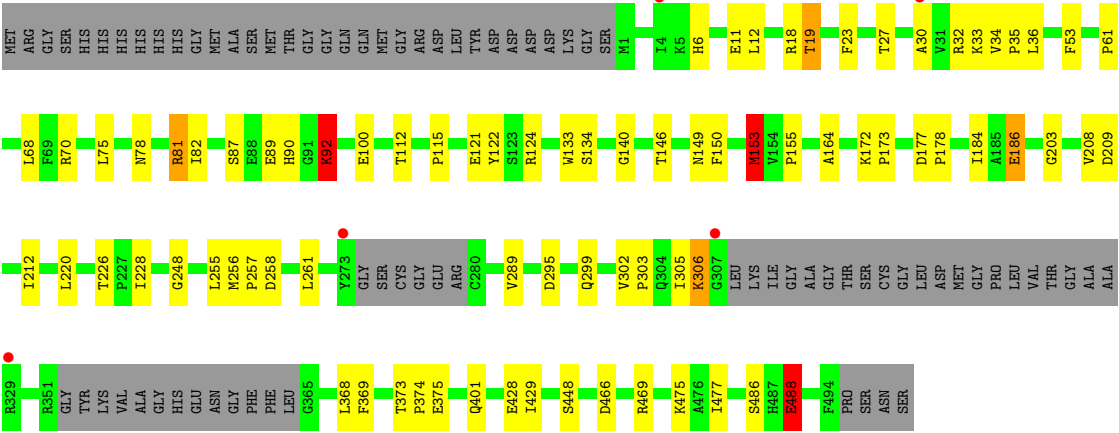


- Molecule 1: Methylmalonate-semialdehyde dehydrogenase





● Molecule 1: Methylmalonate-semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.89Å 156.63Å 192.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.00 – 2.95 49.00 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.00-2.95) 99.9 (49.00-2.95)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.96Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.212 , 0.249 0.216 , 0.251	Depositor DCC
R_{free} test set	3714 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	69.2	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 62.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20975	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% 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: ADP

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	0.77	2/3542 (0.1%)	0.93	2/4810 (0.0%)
1	B	0.77	1/3600 (0.0%)	0.93	4/4889 (0.1%)
1	C	0.85	2/3525 (0.1%)	0.97	6/4787 (0.1%)
1	D	0.75	0/3610	0.96	7/4900 (0.1%)
1	E	0.74	1/3479 (0.0%)	0.91	3/4730 (0.1%)
1	F	0.81	3/3501 (0.1%)	0.96	7/4755 (0.1%)
All	All	0.78	9/21257 (0.0%)	0.94	29/28871 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	D	0	1
1	E	0	1
All	All	0	4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	281	MET	C-O	-6.13	1.16	1.23
1	E	16	THR	C-O	6.09	1.30	1.24
1	B	382	GLU	C-O	-6.05	1.16	1.23
1	A	230	GLU	C-O	-5.58	1.17	1.24
1	A	305	ILE	C-O	-5.52	1.17	1.24
1	F	306	LYS	CA-CB	5.45	1.59	1.52
1	F	19	THR	N-CA	5.45	1.53	1.46
1	C	271	ALA	CA-C	5.15	1.59	1.52
1	F	258	ASP	CG-OD1	-5.12	1.15	1.25

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	454	ARG	CG-CD-NE	9.45	132.80	112.00
1	C	24	ASN	CA-C-N	8.28	130.19	119.84
1	C	24	ASN	C-N-CA	8.28	130.19	119.84
1	D	64	ARG	CG-CD-NE	7.74	129.02	112.00
1	D	24	ASN	N-CA-C	-7.47	96.41	109.48
1	D	244	GLN	CA-CB-CG	7.17	128.44	114.10
1	F	81	ARG	CG-CD-NE	6.91	127.20	112.00
1	F	92	LYS	CA-CB-CG	6.89	127.88	114.10
1	F	488	GLU	CA-CB-CG	6.36	126.81	114.10
1	D	116	GLU	CB-CA-C	-6.03	101.41	110.88
1	F	401	GLN	CB-CG-CD	5.94	122.70	112.60
1	D	129	ASN	N-CA-CB	5.92	120.66	111.84
1	C	28	GLY	N-CA-C	5.84	127.01	113.18
1	F	401	GLN	CA-CB-CG	-5.75	102.61	114.10
1	B	384	PHE	CB-CA-C	5.67	119.09	109.84
1	C	366	GLY	N-CA-C	-5.63	99.84	113.18
1	C	271	ALA	CA-C-O	5.59	128.50	120.51
1	F	153	MET	CB-CG-SD	5.44	129.02	112.70
1	D	175	GLU	O-C-N	-5.42	116.38	122.12
1	F	70	ARG	CB-CG-CD	5.38	123.68	111.30
1	A	230	GLU	CA-C-O	-5.37	114.86	120.55
1	A	117	ILE	CB-CG1-CD1	5.31	124.95	113.80
1	D	494	PHE	CB-CA-C	5.20	119.97	110.10
1	E	382	GLU	CB-CG-CD	5.17	121.39	112.60
1	E	230	GLU	CB-CA-C	5.15	120.16	110.63
1	E	382	GLU	CA-CB-CG	5.09	124.28	114.10
1	C	305	ILE	CA-C-O	5.05	127.10	120.78
1	B	239	ARG	CB-CG-CD	5.05	122.92	111.30
1	B	384	PHE	N-CA-CB	5.04	117.76	109.95

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	230	GLU	Mainchain
1	A	305	ILE	Mainchain
1	D	175	GLU	Mainchain
1	E	260	ASP	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3471	0	3443	56	0
1	B	3525	0	3483	64	0
1	C	3455	0	3410	62	0
1	D	3535	0	3490	69	0
1	E	3408	0	3347	54	0
1	F	3430	0	3405	53	0
2	A	27	0	12	0	0
2	B	27	0	12	1	0
2	D	27	0	12	1	0
2	E	27	0	12	1	0
2	F	27	0	12	1	0
3	A	3	0	0	0	0
3	B	5	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
3	E	1	0	0	0	0
3	F	2	0	0	0	0
All	All	20975	0	20638	349	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 8.

All (349) 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:24:ASN:ND2	1:A:88:GLU:O	1.88	1.06
1:C:46:ILE:CG2	1:C:216:GLU:HG3	1.99	0.92
1:E:269:MET:HE1	1:E:305:ILE:HG13	1.54	0.87
1:C:23:PHE:O	1:C:25:PRO:HD3	1.74	0.86
1:B:152:ALA:O	1:B:156:LEU:HD12	1.79	0.83
1:F:19:THR:HG22	1:F:33:LYS:HG2	1.60	0.83
1:A:209:ASP:HA	1:A:212:ILE:HD12	1.61	0.83
1:E:213:GLU:HA	1:E:239:ARG:NH1	1.93	0.83
1:B:211:LEU:O	1:B:241:LYS:NZ	2.12	0.82
1:B:338:THR:HG22	1:B:342:GLN:HE21	1.43	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:ILE:HG22	1:D:13:ILE:HG12	1.62	0.81
1:C:226:THR:O	1:C:230:GLU:HG3	1.82	0.80
1:B:213:GLU:HG2	1:B:239:ARG:HH22	1.45	0.79
1:A:39:ARG:NH2	1:A:213:GLU:OE1	2.15	0.79
1:E:8:ILE:CD1	1:E:44:GLN:HB3	2.13	0.79
1:D:175:GLU:O	1:D:178:PRO:HD3	1.85	0.76
1:C:46:ILE:HG21	1:C:216:GLU:HG3	1.69	0.75
1:A:208:VAL:HG12	1:A:212:ILE:HD11	1.68	0.75
1:D:175:GLU:O	1:D:178:PRO:HG3	1.86	0.75
1:B:221:SER:OG	1:B:244:GLN:NE2	2.20	0.74
1:E:205:LYS:O	1:E:209:ASP:OD1	2.07	0.72
1:B:269:MET:HE1	1:B:301:LEU:HD23	1.72	0.72
1:A:475:LYS:NZ	1:F:429:ILE:O	2.22	0.72
1:D:8:ILE:HD11	1:D:45:ALA:N	2.05	0.71
1:D:172:LYS:NZ	2:D:501:ADP:O3'	2.24	0.71
1:B:338:THR:HG22	1:B:342:GLN:NE2	2.06	0.70
1:E:119:LYS:NZ	1:F:122:TYR:O	2.23	0.70
1:A:255:LEU:HD11	1:A:297:LEU:HD22	1.74	0.69
1:C:147:PRO:HD3	1:C:223:VAL:HG23	1.74	0.69
1:D:231:TYR:CE1	1:D:235:GLU:OE1	2.45	0.69
1:A:126:VAL:HG21	1:A:132:ALA:HB3	1.75	0.68
1:F:19:THR:CG2	1:F:33:LYS:HG2	2.24	0.68
1:C:23:PHE:C	1:C:25:PRO:HD3	2.21	0.66
1:C:46:ILE:HG21	1:C:216:GLU:CG	2.26	0.66
1:D:429:ILE:O	1:E:475:LYS:NZ	2.29	0.66
1:B:213:GLU:HG2	1:B:239:ARG:NH2	2.12	0.65
1:F:90:HIS:CD2	1:F:92:LYS:HE2	2.32	0.65
1:E:8:ILE:HD11	1:E:44:GLN:C	2.21	0.65
1:D:76:GLU:O	1:D:79:GLU:OE1	2.15	0.64
1:D:475:LYS:NZ	1:E:429:ILE:O	2.30	0.64
1:E:212:ILE:O	1:E:241:LYS:NZ	2.26	0.63
1:F:19:THR:HB	1:F:33:LYS:HE3	1.79	0.63
1:A:80:GLU:O	1:A:84:LYS:HD3	1.99	0.63
1:C:42:MET:SD	1:C:201:VAL:HG11	2.38	0.63
1:B:153:MET:HE1	1:B:456:LEU:HD11	1.81	0.62
1:C:233:TYR:O	1:C:237:THR:HG23	1.98	0.62
1:D:78:ASN:O	1:D:82:ILE:HG13	1.99	0.62
1:D:348:VAL:HG12	1:D:368:LEU:HB3	1.81	0.62
1:F:466:ASP:OD1	1:F:469:ARG:NH1	2.32	0.62
1:B:153:MET:CE	1:B:456:LEU:HD11	2.30	0.62
1:C:201:VAL:O	1:C:201:VAL:HG13	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:GLN:OE1	1:D:473:LYS:NZ	2.32	0.62
1:E:228:ILE:HD12	2:E:501:ADP:O1A	2.00	0.61
1:D:175:GLU:O	1:D:178:PRO:CD	2.49	0.61
1:C:209:ASP:HA	1:C:212:ILE:HG22	1.82	0.61
1:D:257:PRO:HD3	1:D:289:VAL:HG13	1.83	0.61
1:A:24:ASN:CG	1:A:88:GLU:O	2.44	0.60
1:D:8:ILE:HD11	1:D:44:GLN:CB	2.31	0.60
1:F:257:PRO:HD3	1:F:289:VAL:HG13	1.82	0.60
1:C:76:GLU:O	1:C:79:GLU:OE1	2.20	0.59
1:A:215:PRO:HD2	1:A:216:GLU:OE2	2.02	0.59
1:C:78:ASN:O	1:C:82:ILE:HG13	2.01	0.59
1:C:147:PRO:HD3	1:C:223:VAL:CG2	2.32	0.59
1:C:328:ALA:HA	1:C:331:LYS:CG	2.33	0.59
1:B:156:LEU:HD11	1:B:184:ILE:HG21	1.83	0.59
1:D:175:GLU:O	1:D:178:PRO:CG	2.51	0.59
1:B:393:VAL:HG22	1:B:398:GLU:HB3	1.84	0.58
1:F:75:LEU:HD21	1:F:184:ILE:HG23	1.85	0.58
1:B:454:ARG:HG3	1:B:454:ARG:HH11	1.68	0.58
1:E:147:PRO:HD3	1:E:223:VAL:HG13	1.85	0.58
1:B:262:ASP:OD1	1:B:300:LYS:NZ	2.36	0.58
1:F:256:MET:HA	1:F:289:VAL:CG1	2.34	0.58
1:E:74:LEU:HB3	1:E:187:LEU:HD23	1.86	0.57
1:D:256:MET:HA	1:D:289:VAL:CG1	2.34	0.57
1:D:24:ASN:HD22	1:D:25:PRO:CD	2.18	0.57
1:D:64:ARG:NH1	1:D:164:ALA:O	2.37	0.57
1:F:6:HIS:O	1:F:12:LEU:HD23	2.04	0.57
1:E:442:VAL:HG23	1:E:442:VAL:O	2.04	0.57
1:B:75:LEU:HD21	1:B:184:ILE:HG23	1.87	0.57
1:D:213:GLU:HG2	1:D:239:ARG:HH12	1.70	0.57
1:C:388:LEU:HD23	1:C:389:CYS:N	2.20	0.56
1:D:183:LEU:O	1:D:187:LEU:HD22	2.04	0.56
1:C:75:LEU:HD21	1:C:184:ILE:HG23	1.87	0.56
1:D:213:GLU:HA	1:D:239:ARG:HH11	1.70	0.56
1:E:23:PHE:HA	1:E:30:ALA:HA	1.87	0.56
1:A:235:GLU:OE1	1:A:238:LYS:NZ	2.30	0.56
1:A:262:ASP:OD1	1:A:300:LYS:NZ	2.35	0.56
1:B:393:VAL:CG2	1:B:398:GLU:HB3	2.36	0.55
1:D:75:LEU:HD21	1:D:184:ILE:HG23	1.89	0.55
1:E:90:HIS:CD2	1:E:92:LYS:HE2	2.41	0.55
1:C:46:ILE:CG2	1:C:216:GLU:CG	2.80	0.55
1:E:6:HIS:HB2	1:E:13:ILE:HG12	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:8:ILE:HG12	1:E:9:GLY:N	2.22	0.55
1:A:255:LEU:CD1	1:A:297:LEU:HD22	2.36	0.54
1:B:72:LYS:HE3	1:B:109:GLU:HB2	1.89	0.54
1:B:146:THR:HG23	1:B:147:PRO:HD2	1.88	0.54
1:C:116:GLU:OE2	1:C:119:LYS:NZ	2.39	0.54
1:D:8:ILE:CG2	1:D:13:ILE:HG12	2.37	0.54
1:A:146:THR:HG23	1:A:147:PRO:HD2	1.89	0.54
1:B:31:VAL:HG12	1:B:32:ARG:HG3	1.90	0.54
1:C:147:PRO:HG3	1:C:223:VAL:HG23	1.90	0.54
1:C:328:ALA:HA	1:C:331:LYS:HG2	1.89	0.53
1:B:353:TYR:O	1:B:363:PHE:CE1	2.61	0.53
1:F:27:THR:O	1:F:27:THR:HG22	2.08	0.53
1:B:61:PRO:HB3	1:B:115:PRO:HB3	1.91	0.53
1:E:93:THR:HG23	1:E:96:ASP:HB2	1.91	0.53
1:E:213:GLU:HA	1:E:239:ARG:HH11	1.73	0.53
1:E:154:VAL:HG11	1:E:223:VAL:CG1	2.38	0.53
1:A:427:ASP:OD1	1:E:124:ARG:NH2	2.42	0.53
1:B:124[B]:ARG:HH11	1:B:124[B]:ARG:HB2	1.74	0.52
1:D:146:THR:HG23	1:D:147:PRO:HD2	1.91	0.52
1:B:382:GLU:CD	1:B:383:ILE:H	2.17	0.52
1:C:83:VAL:HG23	1:C:97:ALA:HB3	1.91	0.52
1:C:23:PHE:HA	1:C:30:ALA:HA	1.91	0.52
1:E:141:VAL:HG13	1:E:217:VAL:HA	1.91	0.52
1:C:153:MET:HE1	1:C:456:LEU:HG	1.91	0.52
1:B:429:ILE:O	1:C:475:LYS:NZ	2.42	0.52
1:A:118:LEU:HD23	1:A:118:LEU:O	2.10	0.52
1:A:141:VAL:HG13	1:A:217:VAL:HA	1.91	0.52
1:F:295:ASP:O	1:F:299:GLN:HG2	2.10	0.52
1:A:302:VAL:HA	1:A:305:ILE:HD12	1.93	0.51
1:E:301:LEU:O	1:E:305:ILE:HD12	2.11	0.51
1:A:78:ASN:O	1:A:82:ILE:HG13	2.09	0.51
1:C:148:PHE:CD1	1:C:148:PHE:N	2.76	0.51
1:E:212:ILE:C	1:E:241:LYS:NZ	2.69	0.51
1:F:23:PHE:HA	1:F:30:ALA:HA	1.92	0.51
1:C:146:THR:HG21	1:C:155:PRO:HG3	1.92	0.51
1:D:8:ILE:HD11	1:D:44:GLN:HB3	1.90	0.51
1:E:375:GLU:OE1	1:E:375:GLU:N	2.43	0.51
1:A:68:LEU:HB3	1:A:112:THR:HG22	1.93	0.51
1:B:484:ARG:NH1	1:B:491:GLN:O	2.44	0.51
1:D:141:VAL:HG13	1:D:217:VAL:HA	1.93	0.51
1:F:146:THR:HG21	1:F:155:PRO:HG3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:172:LYS:NZ	1:F:173:PRO:O	2.44	0.51
1:D:175:GLU:OE1	1:D:175:GLU:N	2.25	0.50
1:D:255:LEU:O	1:D:289:VAL:HG12	2.10	0.50
1:E:68:LEU:HB3	1:E:112:THR:HG22	1.93	0.50
1:A:235:GLU:HG3	1:A:238:LYS:HE2	1.94	0.50
1:E:146:THR:HG21	1:E:155:PRO:HG3	1.93	0.50
1:A:61:PRO:HB3	1:A:115:PRO:HB3	1.93	0.50
1:F:100:GLU:OE1	1:F:150:PHE:HA	2.12	0.50
1:D:61:PRO:HB3	1:D:115:PRO:HB3	1.94	0.50
1:F:226:THR:OG1	1:F:248:GLY:O	2.29	0.50
1:A:434:VAL:CG1	1:F:477:ILE:HG12	2.42	0.50
1:D:208:VAL:HG12	1:D:212:ILE:HD11	1.94	0.50
1:A:90:HIS:ND1	1:A:92:LYS:HD3	2.27	0.50
1:B:172:LYS:NZ	2:B:501:ADP:O2'	2.32	0.50
1:F:208:VAL:HG12	1:F:212:ILE:HD11	1.94	0.50
1:F:255:LEU:O	1:F:289:VAL:HG12	2.10	0.50
1:D:354:LYS:HB2	1:D:363:PHE:CD1	2.47	0.50
1:F:12:LEU:HD11	1:F:186:GLU:CD	2.37	0.50
1:F:68:LEU:HB3	1:F:112:THR:HG22	1.92	0.50
1:A:208:VAL:O	1:A:212:ILE:HD12	2.12	0.49
1:C:61:PRO:HB3	1:C:115:PRO:HB3	1.94	0.49
1:D:24:ASN:HD22	1:D:25:PRO:HD2	1.76	0.49
1:B:141:VAL:HG13	1:B:217:VAL:HA	1.93	0.49
1:D:226:THR:OG1	1:D:248:GLY:O	2.30	0.49
1:E:226:THR:OG1	1:E:248:GLY:O	2.30	0.49
1:A:235:GLU:HA	1:A:238:LYS:HG2	1.95	0.49
1:B:226:THR:OG1	1:B:248:GLY:O	2.30	0.49
1:B:208:VAL:HG12	1:B:212:ILE:HD11	1.94	0.49
1:A:226:THR:OG1	1:A:248:GLY:O	2.30	0.49
1:A:57:ARG:HH22	1:F:428:GLU:HG2	1.77	0.49
1:D:68:LEU:HB3	1:D:112:THR:HG22	1.93	0.49
1:F:61:PRO:HB3	1:F:115:PRO:HB3	1.95	0.49
1:B:72:LYS:HZ2	1:B:109:GLU:HG3	1.76	0.49
1:C:209:ASP:O	1:C:212:ILE:HG22	2.13	0.49
1:A:373:THR:HG22	1:A:376:MET:HE3	1.95	0.49
1:B:68:LEU:HB3	1:B:112:THR:HG22	1.94	0.49
1:C:226:THR:OG1	1:C:248:GLY:O	2.32	0.48
1:D:24:ASN:HD22	1:D:25:PRO:N	2.11	0.48
1:A:209:ASP:HA	1:A:212:ILE:CD1	2.40	0.48
1:C:328:ALA:HA	1:C:331:LYS:HG3	1.95	0.48
1:D:22:VAL:HG13	1:D:34:VAL:HG13	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:25:PRO:HG2	1:E:90:HIS:C	2.39	0.48
1:B:93:THR:HG22	1:B:95:GLU:HB2	1.95	0.48
1:B:269:MET:HE1	1:B:301:LEU:CD2	2.43	0.48
1:D:81:ARG:O	1:D:85:LEU:HD12	2.13	0.48
1:F:18:ARG:HG2	1:F:36:LEU:HD22	1.95	0.48
1:C:114:ALA:O	1:C:118:LEU:HD13	2.14	0.48
1:D:42:MET:HE3	1:D:201:VAL:HG21	1.96	0.48
1:F:19:THR:HG23	1:F:34:VAL:O	2.14	0.47
1:B:36:LEU:HA	1:B:203:GLY:HA2	1.96	0.47
1:B:23:PHE:HA	1:B:30:ALA:HA	1.96	0.47
1:A:149:ASN:OD1	1:A:149:ASN:N	2.48	0.47
1:D:93:THR:HG23	1:D:96:ASP:HB2	1.96	0.47
1:B:353:TYR:HE1	1:B:362:PHE:CE1	2.33	0.47
1:C:153:MET:HE1	1:C:456:LEU:CG	2.45	0.47
1:C:177:ASP:N	1:C:178:PRO:HD3	2.30	0.47
1:B:250:LYS:HE2	1:B:382:GLU:OE1	2.14	0.47
1:C:69:PHE:CD1	1:C:69:PHE:O	2.68	0.47
1:D:183:LEU:O	1:D:187:LEU:CD2	2.62	0.47
1:E:36:LEU:HA	1:E:203:GLY:HA2	1.96	0.47
1:D:149:ASN:OD1	1:D:149:ASN:N	2.48	0.47
1:B:4:ILE:HD11	1:B:34:VAL:HB	1.96	0.47
1:B:177:ASP:N	1:B:177:ASP:OD1	2.47	0.47
1:C:36:LEU:HA	1:C:203:GLY:HA2	1.96	0.47
1:E:172:LYS:NZ	1:E:173:PRO:O	2.44	0.47
1:A:209:ASP:CA	1:A:212:ILE:HD12	2.38	0.47
1:A:213:GLU:HG2	1:A:239:ARG:NH1	2.30	0.47
1:E:8:ILE:HD12	1:E:44:GLN:HB3	1.92	0.47
1:A:181:THR:O	1:A:184:ILE:HG13	2.15	0.46
1:F:87:SER:OG	1:F:92:LYS:O	2.27	0.46
1:F:177:ASP:N	1:F:178:PRO:HD3	2.30	0.46
1:A:177:ASP:N	1:A:178:PRO:HD3	2.30	0.46
1:F:121:GLU:N	1:F:121:GLU:OE1	2.47	0.46
1:A:292:GLN:CD	1:A:292:GLN:H	2.24	0.46
1:B:87:SER:OG	1:B:92:LYS:O	2.27	0.46
1:C:147:PRO:CD	1:C:223:VAL:HG23	2.43	0.46
1:C:158:MET:HE2	1:C:221:SER:OG	2.14	0.46
1:C:177:ASP:N	1:C:177:ASP:OD1	2.47	0.46
1:D:8:ILE:CD1	1:D:44:GLN:HB3	2.45	0.46
1:D:332:VAL:HG11	1:D:364:LEU:HD22	1.98	0.46
1:F:149:ASN:OD1	1:F:149:ASN:N	2.48	0.46
1:B:149:ASN:OD1	1:B:149:ASN:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ASP:O	1:B:213:GLU:HG3	2.16	0.46
1:B:382:GLU:OE1	1:B:382:GLU:HA	2.16	0.46
1:E:177:ASP:N	1:E:178:PRO:HD3	2.31	0.46
1:F:32:ARG:NH1	1:F:89:GLU:OE2	2.49	0.46
1:F:36:LEU:HA	1:F:203:GLY:HA2	1.97	0.46
1:D:177:ASP:N	1:D:178:PRO:HD3	2.31	0.46
1:E:8:ILE:CD1	1:E:44:GLN:C	2.89	0.46
1:F:177:ASP:N	1:F:177:ASP:OD1	2.48	0.46
1:B:68:LEU:HD11	1:B:164:ALA:HB2	1.96	0.46
1:C:90:HIS:CE1	1:C:92:LYS:HD3	2.51	0.46
1:D:83:VAL:HG11	1:D:94:ILE:HG23	1.98	0.46
1:B:78:ASN:O	1:B:82:ILE:HG13	2.16	0.45
1:B:177:ASP:N	1:B:178:PRO:HD3	2.31	0.45
1:A:36:LEU:HA	1:A:203:GLY:HA2	1.97	0.45
1:B:83:VAL:HG11	1:B:94:ILE:HG23	1.97	0.45
1:D:100:GLU:OE2	1:D:150:PHE:HA	2.17	0.45
1:D:177:ASP:N	1:D:177:ASP:OD1	2.48	0.45
1:E:213:GLU:HA	1:E:239:ARG:HH12	1.74	0.45
1:C:100:GLU:HG3	1:C:153:MET:HB3	1.98	0.45
1:E:149:ASN:OD1	1:E:149:ASN:N	2.48	0.45
1:E:177:ASP:N	1:E:177:ASP:OD1	2.48	0.45
1:A:27:THR:C	1:A:29:GLU:H	2.25	0.45
1:B:100:GLU:OE2	1:B:150:PHE:HA	2.17	0.45
1:C:149:ASN:OD1	1:C:149:ASN:N	2.48	0.45
1:D:332:VAL:CG1	1:D:364:LEU:HD22	2.47	0.45
1:D:68:LEU:HD11	1:D:164:ALA:HB2	1.99	0.45
1:C:306:LYS:HE2	1:C:306:LYS:HA	1.98	0.45
1:C:100:GLU:OE2	1:C:150:PHE:HA	2.16	0.45
1:A:159:TYR:OH	1:A:184:ILE:HD12	2.17	0.44
1:A:333:THR:HG23	1:A:351:ARG:NH1	2.32	0.44
1:E:61:PRO:HB3	1:E:115:PRO:HB3	1.98	0.44
1:F:121:GLU:OE1	1:F:134:SER:OG	2.35	0.44
1:F:228:ILE:HD12	2:F:501:ADP:O2A	2.17	0.44
1:C:117:ILE:HG13	1:C:118:LEU:HD13	1.99	0.44
1:C:154:VAL:HG11	1:C:223:VAL:CG2	2.47	0.44
1:D:8:ILE:HD11	1:D:44:GLN:C	2.42	0.44
1:D:256:MET:SD	1:D:289:VAL:HG11	2.57	0.44
1:A:302:VAL:HB	1:A:303:PRO:HD3	1.99	0.44
1:C:153:MET:HE3	1:C:154:VAL:HG23	1.99	0.44
1:D:36:LEU:HD23	1:D:204:ASP:OD2	2.17	0.44
1:D:137:GLN:OE1	1:D:475:LYS:HE3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:213:GLU:HG2	1:D:239:ARG:NH1	2.32	0.44
1:C:151:PRO:O	1:C:181:THR:HG22	2.16	0.44
1:D:121:GLU:OE2	1:D:121:GLU:N	2.50	0.44
1:C:335:TYR:CE1	1:C:378:ILE:HA	2.53	0.44
1:E:100:GLU:OE2	1:E:150:PHE:HA	2.17	0.44
1:A:68:LEU:HD11	1:A:164:ALA:HB2	1.99	0.44
1:E:133:TRP:CD1	1:F:124:ARG:NH2	2.86	0.44
1:D:83:VAL:HG13	1:D:97:ALA:HB3	2.00	0.44
1:A:100:GLU:OE2	1:A:150:PHE:HA	2.17	0.44
1:E:68:LEU:HD11	1:E:164:ALA:HB2	2.00	0.43
1:F:256:MET:SD	1:F:289:VAL:HG11	2.58	0.43
1:A:87:SER:OG	1:A:92:LYS:O	2.26	0.43
1:A:177:ASP:N	1:A:177:ASP:OD1	2.48	0.43
1:D:36:LEU:HA	1:D:203:GLY:HA2	2.00	0.43
1:F:78:ASN:O	1:F:82:ILE:HG13	2.17	0.43
1:F:368:LEU:C	1:F:368:LEU:HD13	2.44	0.43
1:D:42:MET:HE1	1:D:201:VAL:HB	1.99	0.43
1:E:209:ASP:HA	1:E:212:ILE:HD12	1.99	0.43
1:D:90:HIS:CE1	1:D:92:LYS:HD3	2.52	0.43
1:D:261:LEU:HD13	1:D:293:ILE:HD11	2.01	0.43
1:F:122:TYR:CE1	1:F:133:TRP:CE3	3.06	0.43
1:F:368:LEU:HD13	1:F:369:PHE:N	2.34	0.43
1:B:269:MET:CE	1:B:301:LEU:HD23	2.46	0.43
1:A:217:VAL:HG23	1:A:241:LYS:HE3	2.00	0.43
1:B:8:ILE:HB	1:B:13:ILE:HD12	2.01	0.43
1:B:31:VAL:C	1:B:32:ARG:HG3	2.44	0.43
1:B:302:VAL:HB	1:B:303:PRO:HD3	2.00	0.43
1:F:302:VAL:HB	1:F:303:PRO:HD3	2.00	0.43
1:B:122:TYR:CE2	1:B:133:TRP:CE3	3.07	0.43
1:C:302:VAL:HB	1:C:303:PRO:HD3	2.00	0.43
1:E:122:TYR:CE1	1:E:133:TRP:CE3	3.07	0.43
1:A:122:TYR:CE1	1:A:133:TRP:CE3	3.07	0.43
1:F:68:LEU:HD11	1:F:164:ALA:HB2	2.01	0.43
1:C:46:ILE:HG22	1:C:216:GLU:HG3	1.91	0.42
1:D:8:ILE:CD1	1:D:44:GLN:CB	2.96	0.42
1:D:151:PRO:O	1:D:181:THR:HG22	2.19	0.42
1:A:217:VAL:CG2	1:A:241:LYS:HE3	2.49	0.42
1:F:209:ASP:HA	1:F:212:ILE:HD12	2.01	0.42
1:A:93:THR:OG1	1:A:95:GLU:OE1	2.37	0.42
1:D:302:VAL:HB	1:D:303:PRO:HD3	2.01	0.42
1:D:122:TYR:CE1	1:D:133:TRP:CE3	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:PHE:CZ	1:A:140:GLY:HA2	2.55	0.42
1:C:27:THR:O	1:C:27:THR:OG1	2.29	0.42
1:C:40:GLU:N	1:C:40:GLU:OE1	2.52	0.42
1:E:154:VAL:HG11	1:E:223:VAL:HG13	2.01	0.42
1:B:53:PHE:CZ	1:B:140:GLY:HA2	2.54	0.42
1:E:302:VAL:HB	1:E:303:PRO:HD3	2.00	0.42
1:B:261:LEU:HD13	1:B:293:ILE:HD11	2.02	0.42
1:E:217:VAL:CG2	1:E:241:LYS:HE2	2.50	0.42
1:B:126:VAL:HG21	1:B:132:ALA:HB3	2.01	0.42
1:C:83:VAL:HG23	1:C:97:ALA:CB	2.49	0.42
1:B:335:TYR:CE1	1:B:378:ILE:HA	2.55	0.41
1:D:209:ASP:HA	1:D:212:ILE:HD12	2.02	0.41
1:E:2:THR:HG22	1:E:3:LEU:N	2.35	0.41
1:E:87:SER:OG	1:E:92:LYS:O	2.28	0.41
1:C:217:VAL:HG23	1:C:241:LYS:HE3	2.01	0.41
1:D:93:THR:OG1	1:D:95:GLU:OE1	2.38	0.41
1:C:53:PHE:CZ	1:C:140:GLY:HA2	2.55	0.41
1:C:68:LEU:HD11	1:C:164:ALA:HB2	2.03	0.41
1:F:486:SER:HB2	1:F:488:GLU:OE1	2.20	0.41
1:B:118:LEU:O	1:B:118:LEU:HD23	2.20	0.41
1:A:335:TYR:CE1	1:A:378:ILE:HA	2.56	0.41
1:F:53:PHE:CZ	1:F:140:GLY:HA2	2.55	0.41
1:B:83:VAL:HG13	1:B:97:ALA:HB3	2.02	0.41
1:B:233:TYR:HA	1:B:243:VAL:HG11	2.02	0.41
1:C:217:VAL:CG2	1:C:241:LYS:HE3	2.50	0.41
1:D:233:TYR:HA	1:D:243:VAL:HG11	2.03	0.41
1:E:233:TYR:HA	1:E:243:VAL:HG11	2.03	0.41
1:F:261:LEU:HD23	1:F:261:LEU:HA	1.92	0.41
1:A:2:THR:HG22	1:A:3:LEU:N	2.35	0.41
1:B:364:LEU:HD23	1:B:365:GLY:O	2.20	0.41
1:C:34:VAL:HA	1:C:35:PRO:HD3	1.94	0.41
1:C:122:TYR:CE2	1:C:133:TRP:CE3	3.09	0.41
1:C:328:ALA:N	1:C:331:LYS:HZ3	2.18	0.41
1:A:82:ILE:HD13	1:A:184:ILE:CG2	2.51	0.41
1:A:213:GLU:HG2	1:A:239:ARG:HH12	1.85	0.41
1:A:233:TYR:HA	1:A:243:VAL:HG11	2.03	0.41
1:E:34:VAL:HA	1:E:35:PRO:HD3	1.93	0.41
1:F:100:GLU:HG3	1:F:153:MET:HB3	2.03	0.41
1:F:373:THR:CG2	1:F:374:PRO:HD2	2.50	0.41
1:D:335:TYR:CE1	1:D:378:ILE:HA	2.55	0.41
1:E:335:TYR:CE1	1:E:378:ILE:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:TYR:HA	1:C:243:VAL:HG11	2.02	0.41
1:E:53:PHE:CZ	1:E:140:GLY:HA2	2.56	0.41
1:F:34:VAL:HA	1:F:35:PRO:HD3	1.93	0.41
1:F:305:ILE:O	1:F:306:LYS:HG2	2.21	0.41
1:B:124[A]:ARG:HE	1:B:124[A]:ARG:HB3	1.74	0.40
1:B:152:ALA:O	1:B:156:LEU:CD1	2.59	0.40
1:A:126:VAL:HG21	1:A:132:ALA:CB	2.49	0.40
1:E:117:ILE:O	1:E:117:ILE:HG13	2.21	0.40
1:B:38:ASP:OD1	1:B:38:ASP:N	2.55	0.40
1:E:8:ILE:HG13	1:E:45:ALA:HA	2.03	0.40
1:F:220:LEU:HD12	1:F:220:LEU:HA	1.98	0.40

There are no symmetry-related clashes.

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	452/531 (85%)	432 (96%)	20 (4%)	0	100	100
1	B	458/531 (86%)	438 (96%)	20 (4%)	0	100	100
1	C	453/531 (85%)	430 (95%)	21 (5%)	2 (0%)	30	53
1	D	460/531 (87%)	444 (96%)	16 (4%)	0	100	100
1	E	447/531 (84%)	428 (96%)	19 (4%)	0	100	100
1	F	446/531 (84%)	423 (95%)	23 (5%)	0	100	100
All	All	2716/3186 (85%)	2595 (96%)	119 (4%)	2 (0%)	48	72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	28	GLY
1	C	271	ALA

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	353/416 (85%)	341 (97%)	12 (3%)	32	58
1	B	358/416 (86%)	346 (97%)	12 (3%)	32	58
1	C	349/416 (84%)	337 (97%)	12 (3%)	32	58
1	D	358/416 (86%)	342 (96%)	16 (4%)	24	50
1	E	343/416 (82%)	335 (98%)	8 (2%)	44	67
1	F	350/416 (84%)	341 (97%)	9 (3%)	40	65
All	All	2111/2496 (85%)	2042 (97%)	69 (3%)	34	58

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	21	ASP
1	A	39	ARG
1	A	80	GLU
1	A	101	LEU
1	A	118	LEU
1	A	119	LYS
1	A	149	ASN
1	A	284	SER
1	A	382	GLU
1	A	448	SER
1	A	491	GLN
1	B	4	ILE
1	B	38	ASP
1	B	92	LYS
1	B	124[A]	ARG
1	B	124[B]	ARG
1	B	153	MET
1	B	284	SER
1	B	353	TYR
1	B	375	GLU
1	B	454	ARG

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Mol	Chain	Res	Type
1	B	484	ARG
1	B	488	GLU
1	C	32	ARG
1	C	57	ARG
1	C	66	GLN
1	C	80	GLU
1	C	81	ARG
1	C	101	LEU
1	C	124	ARG
1	C	148	PHE
1	C	153	MET
1	C	179	SER
1	C	216	GLU
1	C	453	LYS
1	D	24	ASN
1	D	57	ARG
1	D	82	ILE
1	D	85	LEU
1	D	116	GLU
1	D	129	ASN
1	D	149	ASN
1	D	244	GLN
1	D	300	LYS
1	D	348	VAL
1	D	375	GLU
1	D	401	GLN
1	D	448	SER
1	D	469	ARG
1	D	475	LYS
1	D	488	GLU
1	E	32	ARG
1	E	63	LYS
1	E	81	ARG
1	E	92	LYS
1	E	187	LEU
1	E	223	VAL
1	E	473	LYS
1	E	474	ARG
1	F	11	GLU
1	F	81	ARG
1	F	92	LYS
1	F	153	MET

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Mol	Chain	Res	Type
1	F	186	GLU
1	F	375	GLU
1	F	448	SER
1	F	475	LYS
1	F	488	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	58	ASN
1	A	66	GLN
1	A	202	HIS
1	A	491	GLN
1	B	129	ASN
1	B	202	HIS
1	B	244	GLN
1	B	292	GLN
1	B	342	GLN
1	B	491	GLN
1	C	73	GLN
1	C	263	ASN
1	C	292	GLN
1	C	304	GLN
1	D	6	HIS
1	D	24	ASN
1	D	44	GLN
1	D	90	HIS
1	D	202	HIS
1	D	360	ASN
1	E	58	ASN
1	E	78	ASN
1	E	189	HIS
1	E	491	GLN
1	F	125	ASN
1	F	129	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	$\# Z > 2$
2	ADP	E	501	-	28,29,29	1.53	5 (17%)	43,45,45	2.12	10 (23%)
2	ADP	F	501	-	28,29,29	1.38	3 (10%)	43,45,45	2.09	12 (27%)
2	ADP	D	501	-	28,29,29	1.47	4 (14%)	43,45,45	1.97	10 (23%)
2	ADP	B	501	-	28,29,29	1.49	5 (17%)	43,45,45	2.16	13 (30%)
2	ADP	A	501	-	28,29,29	1.46	4 (14%)	43,45,45	2.05	12 (27%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	E	501	-	-	3/16/32/32	0/3/3/3
2	ADP	F	501	-	-	9/16/32/32	0/3/3/3
2	ADP	D	501	-	-	7/16/32/32	0/3/3/3
2	ADP	B	501	-	-	5/16/32/32	0/3/3/3
2	ADP	A	501	-	-	9/16/32/32	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	501	ADP	C5-C4	4.67	1.47	1.39
2	D	501	ADP	C5-C4	4.52	1.47	1.39
2	A	501	ADP	C5-C4	4.46	1.47	1.39
2	F	501	ADP	C5-C4	4.44	1.47	1.39
2	B	501	ADP	C5-C4	4.23	1.46	1.39
2	E	501	ADP	PA-O3A	3.31	1.63	1.59
2	F	501	ADP	C5-N7	-3.30	1.33	1.39
2	A	501	ADP	C5-C6	3.27	1.50	1.41
2	D	501	ADP	PA-O3A	2.88	1.62	1.59
2	E	501	ADP	C5-C6	2.86	1.49	1.41
2	B	501	ADP	PA-O3A	2.82	1.62	1.59
2	B	501	ADP	C5-C6	2.75	1.48	1.41
2	A	501	ADP	C4-N9	-2.64	1.32	1.37
2	A	501	ADP	C8-N7	2.57	1.36	1.31
2	E	501	ADP	C5-N7	-2.55	1.34	1.39
2	D	501	ADP	C5-C6	2.41	1.47	1.41
2	B	501	ADP	C5-N7	-2.40	1.34	1.39
2	D	501	ADP	C5-N7	-2.39	1.34	1.39
2	E	501	ADP	C8-N7	2.08	1.35	1.31
2	F	501	ADP	C5-C6	2.05	1.46	1.41
2	B	501	ADP	C8-N7	2.04	1.35	1.31

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	501	ADP	C5-C4-N3	-7.45	116.47	126.72
2	F	501	ADP	C5-C4-N3	-6.73	117.46	126.72
2	B	501	ADP	C5-C4-N3	-6.27	118.08	126.72
2	E	501	ADP	N3-C4-N9	6.11	137.56	127.17
2	F	501	ADP	N3-C4-N9	5.93	137.25	127.17
2	A	501	ADP	C5-C4-N3	-5.71	118.85	126.72
2	D	501	ADP	C5-C4-N3	-5.44	119.22	126.72
2	B	501	ADP	N3-C4-N9	5.30	136.18	127.17
2	D	501	ADP	N3-C4-N9	4.96	135.61	127.17
2	E	501	ADP	C2-N3-C4	4.55	122.95	111.83
2	A	501	ADP	C4-C5-N7	-4.53	105.40	110.58
2	D	501	ADP	N3-C2-N1	-4.39	121.94	128.58
2	A	501	ADP	N3-C4-N9	4.14	134.20	127.17
2	B	501	ADP	C2-N3-C4	4.10	121.84	111.83
2	F	501	ADP	C2-N3-C4	3.95	121.47	111.83
2	D	501	ADP	C2-N3-C4	3.87	121.27	111.83
2	A	501	ADP	C4-N9-C8	3.80	109.72	105.74
2	A	501	ADP	C2-N3-C4	3.72	120.92	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	501	ADP	N3-C2-N1	-3.71	122.97	128.58
2	D	501	ADP	C4-N9-C8	3.62	109.54	105.74
2	B	501	ADP	C4-N9-C8	3.60	109.52	105.74
2	B	501	ADP	C4-C5-N7	-3.58	106.49	110.58
2	E	501	ADP	N3-C2-N1	-3.56	123.19	128.58
2	B	501	ADP	N3-C2-N1	-3.41	123.42	128.58
2	E	501	ADP	C4-C5-N7	-3.28	106.83	110.58
2	A	501	ADP	C5-N7-C8	3.23	108.53	103.45
2	B	501	ADP	C5-N7-C8	3.22	108.51	103.45
2	A	501	ADP	N9-C8-N7	-3.20	109.39	113.94
2	A	501	ADP	C6-C5-N7	3.10	138.06	132.09
2	F	501	ADP	C4-N9-C8	3.00	108.88	105.74
2	E	501	ADP	C5-N7-C8	2.98	108.14	103.45
2	F	501	ADP	C5-N7-C8	2.98	108.13	103.45
2	B	501	ADP	N9-C8-N7	-2.97	109.72	113.94
2	F	501	ADP	C4-C5-N7	-2.96	107.19	110.58
2	B	501	ADP	C5'-C4'-C3'	-2.92	104.68	115.21
2	A	501	ADP	N3-C2-N1	-2.85	124.27	128.58
2	D	501	ADP	C5-N7-C8	2.77	107.80	103.45
2	D	501	ADP	N9-C8-N7	-2.75	110.04	113.94
2	D	501	ADP	C4-C5-N7	-2.74	107.45	110.58
2	E	501	ADP	C3'-C2'-C1'	2.66	106.50	101.46
2	A	501	ADP	C4-N9-C1'	-2.59	120.57	126.63
2	B	501	ADP	O2A-PA-O3A	2.58	114.26	107.27
2	D	501	ADP	C2-N1-C6	2.57	122.95	118.73
2	B	501	ADP	O3'-C3'-C4'	-2.54	103.77	111.08
2	A	501	ADP	O3B-PB-O2B	2.52	117.26	107.80
2	F	501	ADP	N6-C6-N1	2.47	123.87	118.38
2	A	501	ADP	O2A-PA-O1A	2.45	123.83	112.44
2	F	501	ADP	N9-C8-N7	-2.43	110.48	113.94
2	E	501	ADP	C4-N9-C8	2.38	108.23	105.74
2	E	501	ADP	C2'-C3'-C4'	2.23	106.92	102.61
2	E	501	ADP	N9-C8-N7	-2.20	110.82	113.94
2	B	501	ADP	C6-C5-N7	2.19	136.32	132.09
2	F	501	ADP	O3B-PB-O2B	2.15	115.86	107.80
2	F	501	ADP	C5-C6-N6	-2.13	118.02	123.29
2	B	501	ADP	O2A-PA-O1A	2.09	122.16	112.44
2	D	501	ADP	O3A-PB-O1B	-2.05	100.25	111.04
2	F	501	ADP	C3'-C2'-C1'	2.02	105.28	101.46

There are no chirality outliers.

All (33) torsion outliers are listed below:

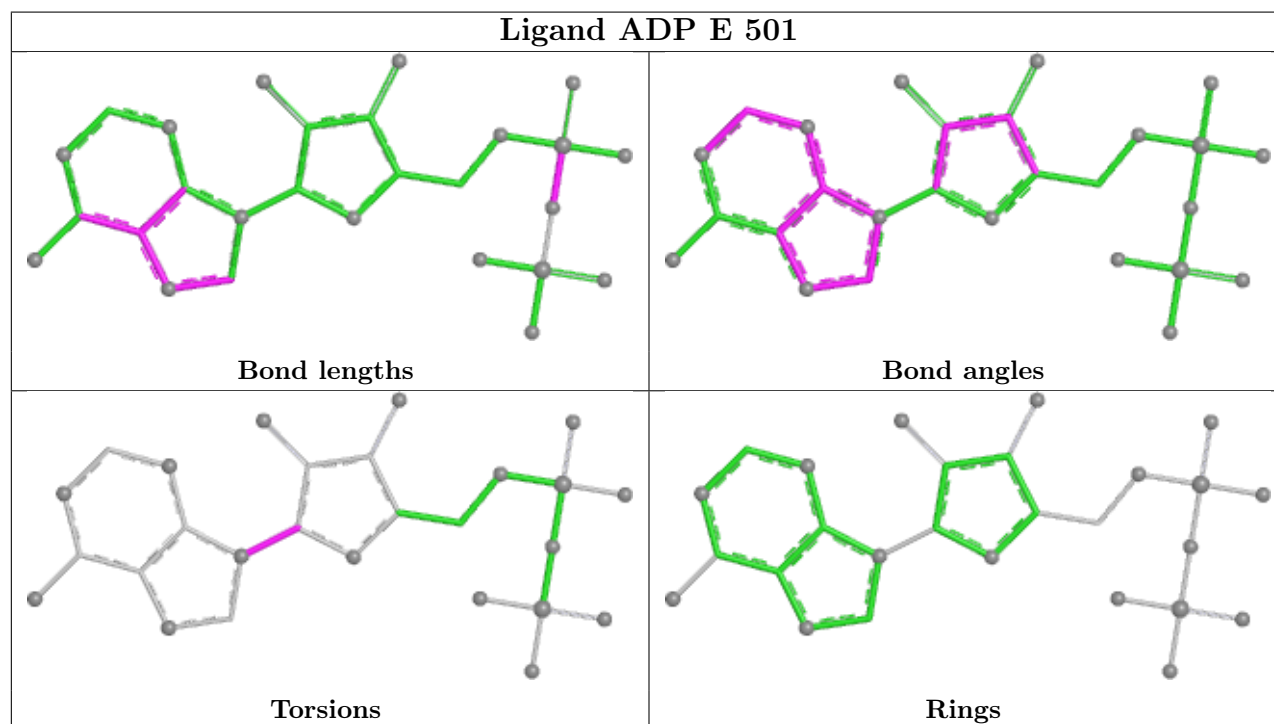
Mol	Chain	Res	Type	Atoms
2	A	501	ADP	PA-O3A-PB-O2B
2	A	501	ADP	O4'-C4'-C5'-O5'
2	B	501	ADP	PA-O3A-PB-O3B
2	D	501	ADP	C5'-O5'-PA-O3A
2	F	501	ADP	C5'-O5'-PA-O1A
2	F	501	ADP	C5'-O5'-PA-O3A
2	F	501	ADP	O4'-C4'-C5'-O5'
2	F	501	ADP	C3'-C4'-C5'-O5'
2	A	501	ADP	C3'-C4'-C5'-O5'
2	D	501	ADP	O4'-C4'-C5'-O5'
2	E	501	ADP	C2'-C1'-N9-C8
2	D	501	ADP	C3'-C4'-C5'-O5'
2	F	501	ADP	PA-O3A-PB-O1B
2	A	501	ADP	C2'-C1'-N9-C8
2	D	501	ADP	C2'-C1'-N9-C8
2	F	501	ADP	C2'-C1'-N9-C8
2	A	501	ADP	C5'-O5'-PA-O1A
2	A	501	ADP	C5'-O5'-PA-O2A
2	A	501	ADP	C5'-O5'-PA-O3A
2	A	501	ADP	C2'-C1'-N9-C4
2	D	501	ADP	C2'-C1'-N9-C4
2	E	501	ADP	C2'-C1'-N9-C4
2	D	501	ADP	O4'-C1'-N9-C8
2	B	501	ADP	PA-O3A-PB-O1B
2	B	501	ADP	PA-O3A-PB-O2B
2	F	501	ADP	PA-O3A-PB-O2B
2	F	501	ADP	PA-O3A-PB-O3B
2	E	501	ADP	O4'-C1'-N9-C8
2	B	501	ADP	C3'-C4'-C5'-O5'
2	D	501	ADP	PB-O3A-PA-O2A
2	A	501	ADP	O4'-C1'-N9-C8
2	F	501	ADP	O4'-C1'-N9-C8
2	B	501	ADP	PB-O3A-PA-O1A

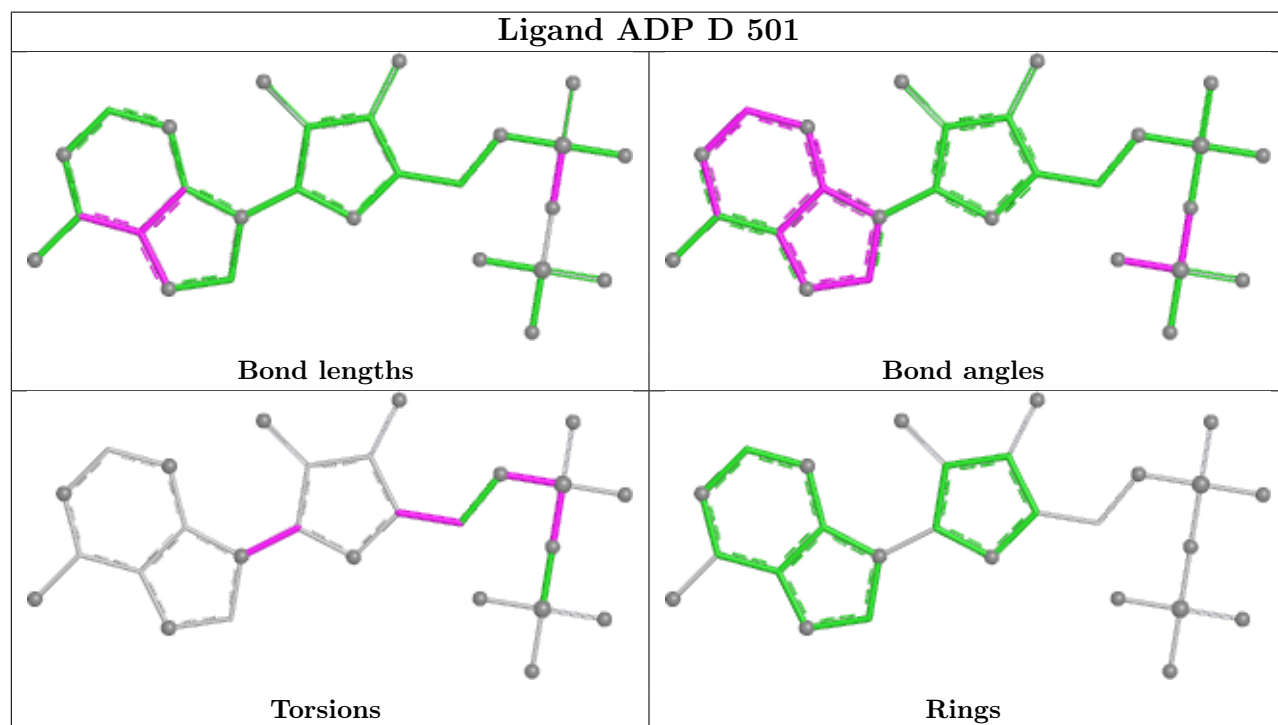
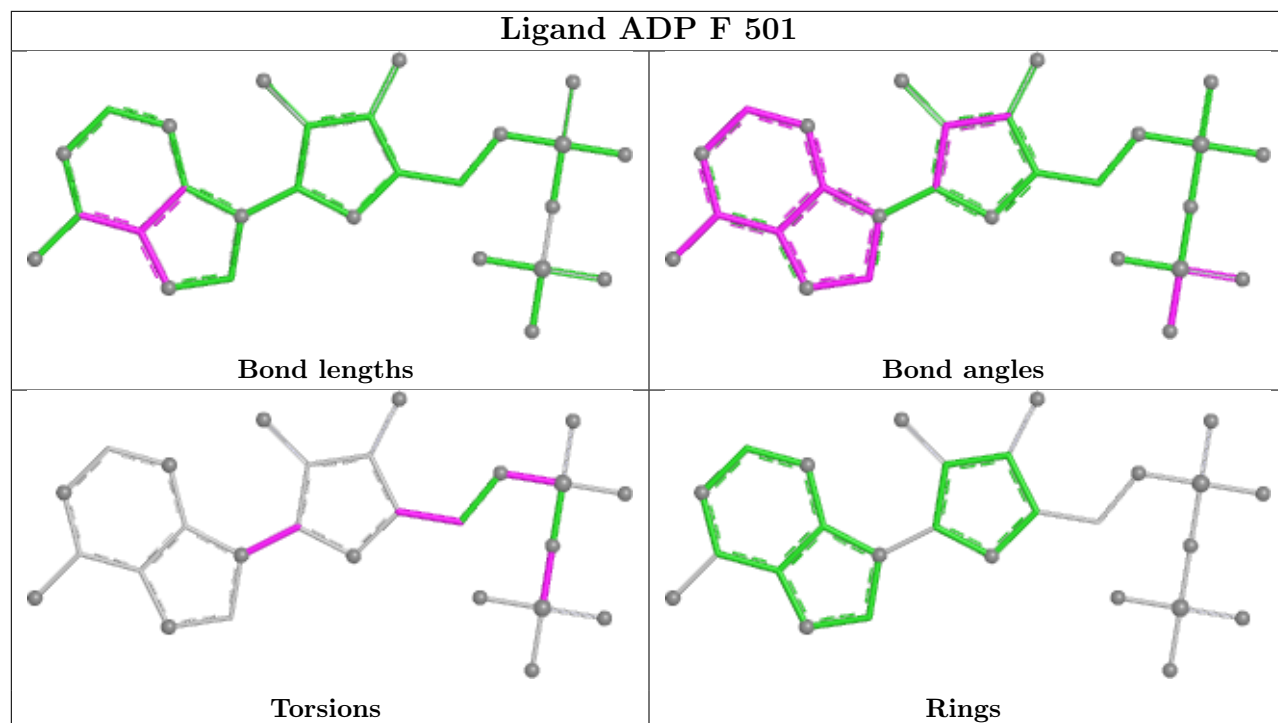
There are no ring outliers.

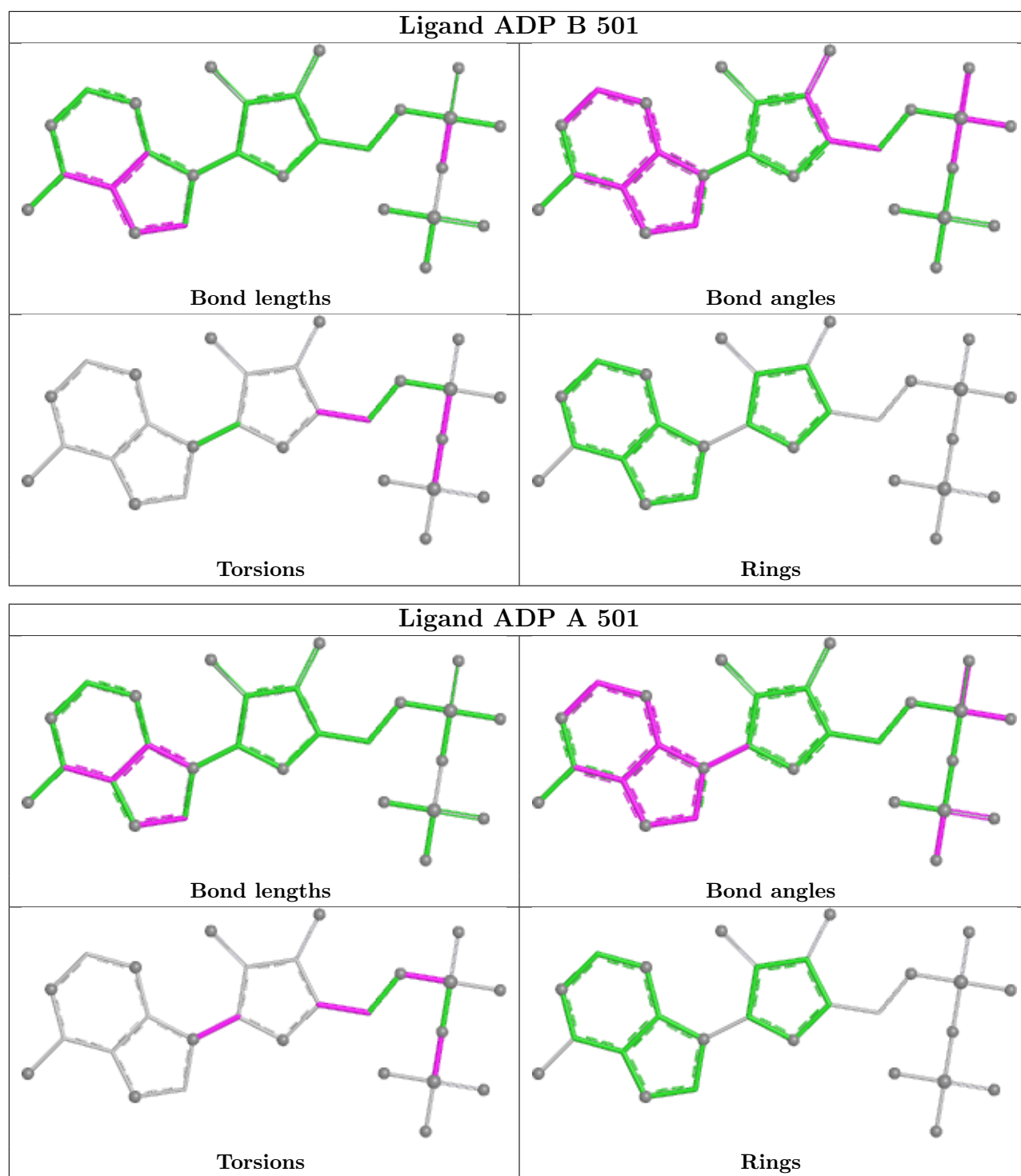
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	501	ADP	1	0
2	F	501	ADP	1	0
2	D	501	ADP	1	0
2	B	501	ADP	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

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	460/531 (86%)	-0.00	7 (1%) 72 65	38, 65, 115, 157	0
1	B	465/531 (87%)	0.02	7 (1%) 72 65	31, 65, 118, 149	1 (0%)
1	C	461/531 (86%)	0.11	16 (3%) 47 40	34, 67, 114, 148	0
1	D	468/531 (88%)	0.05	7 (1%) 72 65	36, 67, 114, 141	0
1	E	455/531 (85%)	0.17	6 (1%) 75 69	39, 73, 126, 157	0
1	F	454/531 (85%)	0.09	5 (1%) 78 73	39, 69, 111, 139	0
All	All	2763/3186 (86%)	0.07	48 (1%) 69 62	31, 68, 118, 157	1 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	364	LEU	3.7
1	C	272	ALA	3.3
1	C	201	VAL	3.2
1	C	26	SER	2.9
1	E	493	ALA	2.8
1	A	96	ASP	2.8
1	C	273	TYR	2.7
1	D	274	GLY	2.7
1	C	328	ALA	2.7
1	C	365	GLY	2.7
1	F	307	GLY	2.6
1	D	494	PHE	2.6
1	B	31	VAL	2.6
1	B	118	LEU	2.6
1	A	329	ARG	2.6
1	C	308	LEU	2.6
1	B	117	ILE	2.6
1	F	4	ILE	2.5
1	F	329	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	355	VAL	2.5
1	C	352	GLY	2.5
1	A	273	TYR	2.4
1	A	327	ALA	2.4
1	B	2	THR	2.3
1	C	274	GLY	2.3
1	E	274	GLY	2.3
1	A	278	GLU	2.3
1	C	271	ALA	2.3
1	A	4	ILE	2.3
1	D	4	ILE	2.3
1	F	273	TYR	2.3
1	C	332	VAL	2.2
1	A	381	GLU	2.2
1	E	133	TRP	2.2
1	C	279	ARG	2.1
1	C	384	PHE	2.1
1	D	299	GLN	2.1
1	B	493	ALA	2.1
1	E	184	ILE	2.1
1	F	30	ALA	2.1
1	B	494	PHE	2.0
1	E	201	VAL	2.0
1	C	307	GLY	2.0
1	D	384	PHE	2.0
1	C	4	ILE	2.0
1	D	27	THR	2.0
1	B	133	TRP	2.0
1	E	273	TYR	2.0

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

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

6.3 Carbohydrates ⓘ

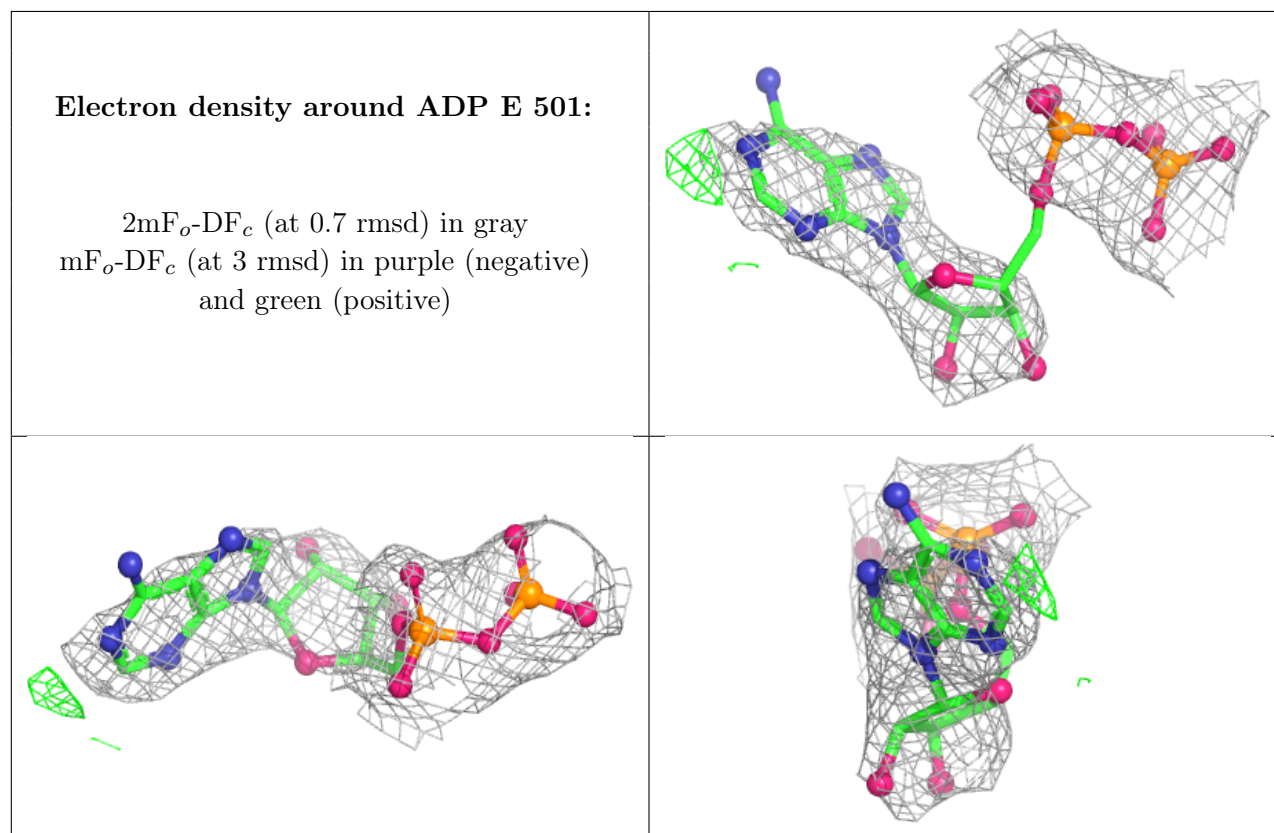
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

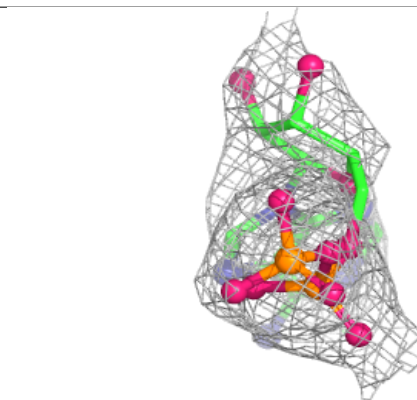
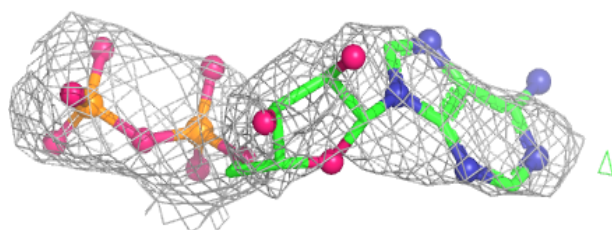
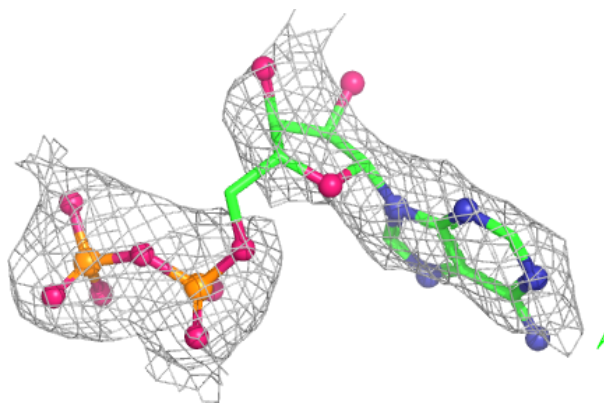
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ADP	E	501	27/27	0.84	0.15	83,95,105,108	27
2	ADP	A	501	27/27	0.87	0.14	76,91,100,105	27
2	ADP	F	501	27/27	0.88	0.13	65,80,89,95	27
2	ADP	D	501	27/27	0.89	0.11	59,67,74,77	27
2	ADP	B	501	27/27	0.91	0.12	61,67,72,76	27

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.

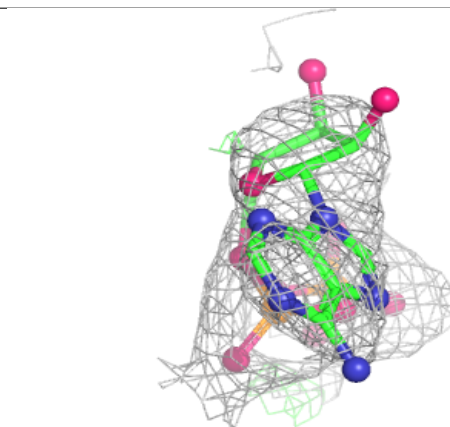
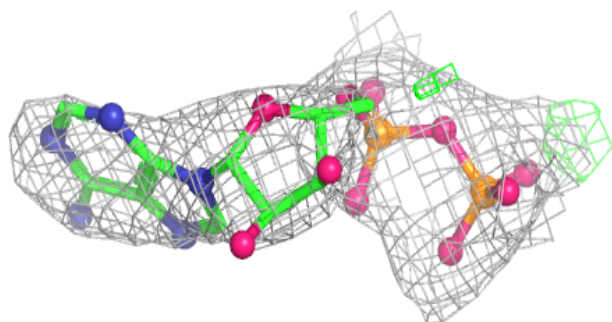
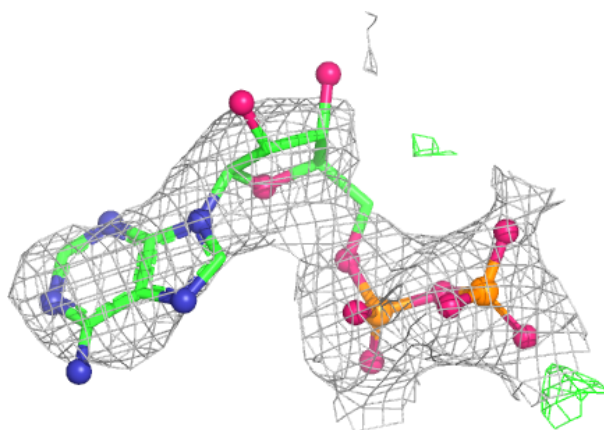


Electron density around ADP A 501:

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

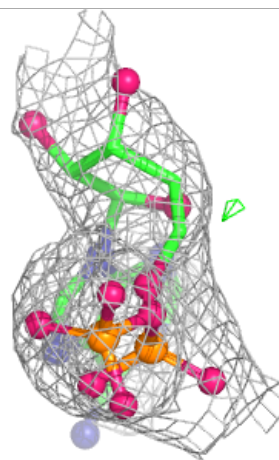
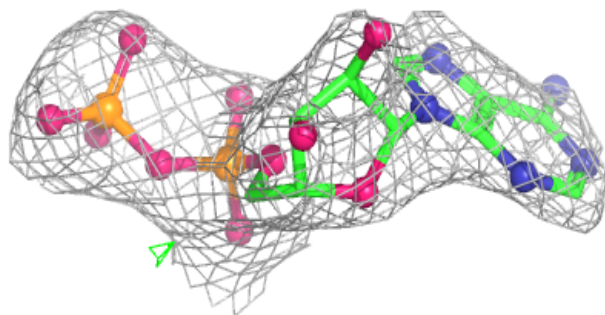
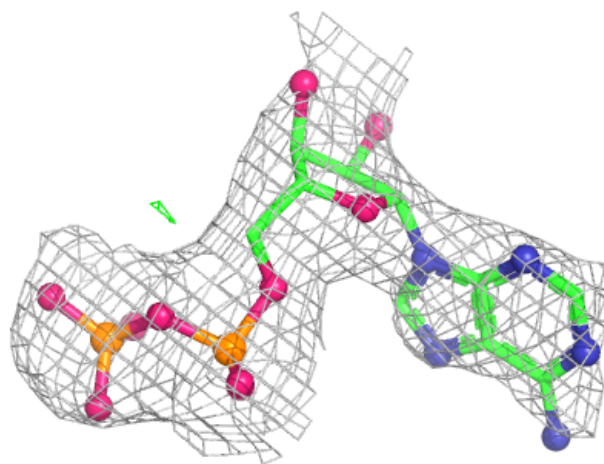
**Electron density around ADP F 501:**

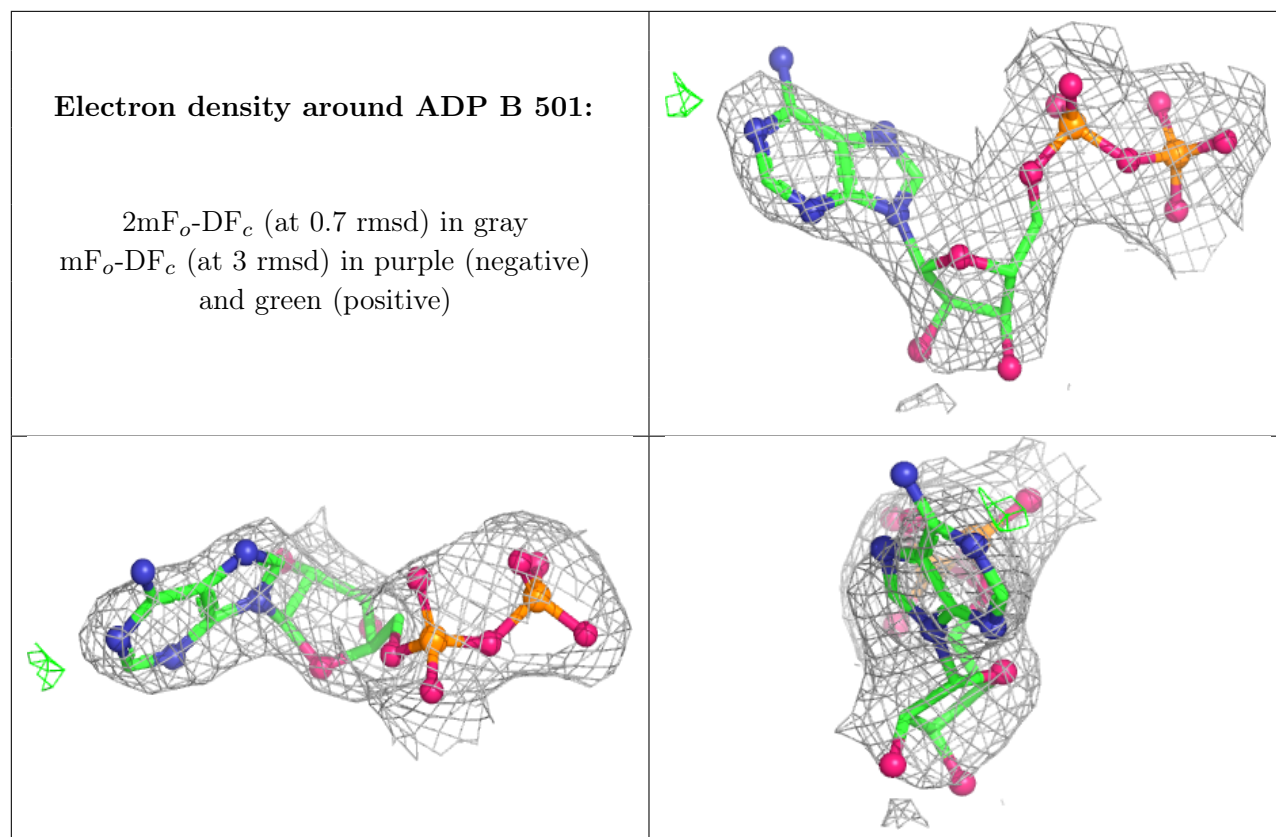
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ADP D 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 [i](#)

There are no such residues in this entry.