



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:11 AM UTC

PDB ID : 6QAO / pdb_00006qao
Title : Structure of human aldehyde dehydrogenase 9A1 in P21 space group
Authors : Morera, S.; Vigouroux, A.
Deposited on : 2018-12-19
Resolution : 2.89 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

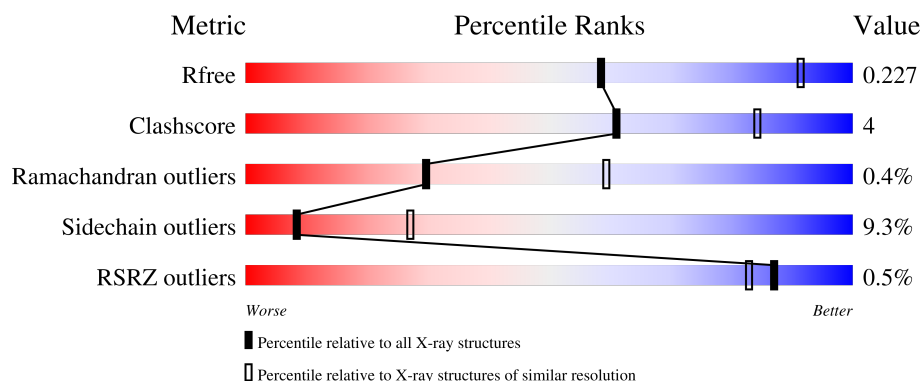
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

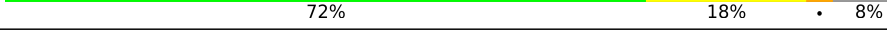
The reported resolution of this entry is 2.89 Å.

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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	508	
1	B	508	
1	C	508	
1	D	508	
1	E	508	

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Mol	Chain	Length	Quality of chain
1	F	508	76% 14% • 7%
1	G	508	73% 16% • 8%
1	H	508	% 76% 13% • 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 28772 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 4-trimethylaminobutyraldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	0	0
			3600	2279	615	678	28			
1	B	469	Total	C	N	O	S	0	0	0
			3596	2277	614	677	28			
1	C	470	Total	C	N	O	S	0	0	0
			3600	2279	615	678	28			
1	D	468	Total	C	N	O	S	0	1	0
			3594	2277	613	675	29			
1	E	465	Total	C	N	O	S	0	0	0
			3567	2258	610	671	28			
1	F	470	Total	C	N	O	S	0	0	0
			3597	2277	615	677	28			
1	G	469	Total	C	N	O	S	0	0	0
			3596	2277	614	677	28			
1	H	468	Total	C	N	O	S	0	0	0
			3589	2273	613	675	28			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP P49189
A	-12	GLY	-	expression tag	UNP P49189
A	-11	SER	-	expression tag	UNP P49189
A	-10	SER	-	expression tag	UNP P49189
A	-9	HIS	-	expression tag	UNP P49189
A	-8	HIS	-	expression tag	UNP P49189
A	-7	HIS	-	expression tag	UNP P49189
A	-6	HIS	-	expression tag	UNP P49189
A	-5	HIS	-	expression tag	UNP P49189
A	-4	HIS	-	expression tag	UNP P49189
A	-3	SER	-	expression tag	UNP P49189
A	-2	GLN	-	expression tag	UNP P49189
A	-1	ASP	-	expression tag	UNP P49189

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	PRO	-	expression tag	UNP P49189
B	-13	MET	-	initiating methionine	UNP P49189
B	-12	GLY	-	expression tag	UNP P49189
B	-11	SER	-	expression tag	UNP P49189
B	-10	SER	-	expression tag	UNP P49189
B	-9	HIS	-	expression tag	UNP P49189
B	-8	HIS	-	expression tag	UNP P49189
B	-7	HIS	-	expression tag	UNP P49189
B	-6	HIS	-	expression tag	UNP P49189
B	-5	HIS	-	expression tag	UNP P49189
B	-4	HIS	-	expression tag	UNP P49189
B	-3	SER	-	expression tag	UNP P49189
B	-2	GLN	-	expression tag	UNP P49189
B	-1	ASP	-	expression tag	UNP P49189
B	0	PRO	-	expression tag	UNP P49189
C	-13	MET	-	initiating methionine	UNP P49189
C	-12	GLY	-	expression tag	UNP P49189
C	-11	SER	-	expression tag	UNP P49189
C	-10	SER	-	expression tag	UNP P49189
C	-9	HIS	-	expression tag	UNP P49189
C	-8	HIS	-	expression tag	UNP P49189
C	-7	HIS	-	expression tag	UNP P49189
C	-6	HIS	-	expression tag	UNP P49189
C	-5	HIS	-	expression tag	UNP P49189
C	-4	HIS	-	expression tag	UNP P49189
C	-3	SER	-	expression tag	UNP P49189
C	-2	GLN	-	expression tag	UNP P49189
C	-1	ASP	-	expression tag	UNP P49189
C	0	PRO	-	expression tag	UNP P49189
D	-13	MET	-	initiating methionine	UNP P49189
D	-12	GLY	-	expression tag	UNP P49189
D	-11	SER	-	expression tag	UNP P49189
D	-10	SER	-	expression tag	UNP P49189
D	-9	HIS	-	expression tag	UNP P49189
D	-8	HIS	-	expression tag	UNP P49189
D	-7	HIS	-	expression tag	UNP P49189
D	-6	HIS	-	expression tag	UNP P49189
D	-5	HIS	-	expression tag	UNP P49189
D	-4	HIS	-	expression tag	UNP P49189
D	-3	SER	-	expression tag	UNP P49189
D	-2	GLN	-	expression tag	UNP P49189
D	-1	ASP	-	expression tag	UNP P49189

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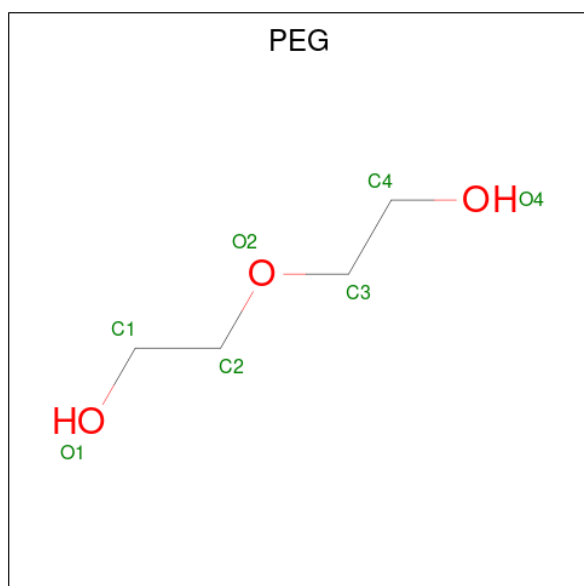
Chain	Residue	Modelled	Actual	Comment	Reference
D	0	PRO	-	expression tag	UNP P49189
E	-13	MET	-	initiating methionine	UNP P49189
E	-12	GLY	-	expression tag	UNP P49189
E	-11	SER	-	expression tag	UNP P49189
E	-10	SER	-	expression tag	UNP P49189
E	-9	HIS	-	expression tag	UNP P49189
E	-8	HIS	-	expression tag	UNP P49189
E	-7	HIS	-	expression tag	UNP P49189
E	-6	HIS	-	expression tag	UNP P49189
E	-5	HIS	-	expression tag	UNP P49189
E	-4	HIS	-	expression tag	UNP P49189
E	-3	SER	-	expression tag	UNP P49189
E	-2	GLN	-	expression tag	UNP P49189
E	-1	ASP	-	expression tag	UNP P49189
E	0	PRO	-	expression tag	UNP P49189
F	-13	MET	-	initiating methionine	UNP P49189
F	-12	GLY	-	expression tag	UNP P49189
F	-11	SER	-	expression tag	UNP P49189
F	-10	SER	-	expression tag	UNP P49189
F	-9	HIS	-	expression tag	UNP P49189
F	-8	HIS	-	expression tag	UNP P49189
F	-7	HIS	-	expression tag	UNP P49189
F	-6	HIS	-	expression tag	UNP P49189
F	-5	HIS	-	expression tag	UNP P49189
F	-4	HIS	-	expression tag	UNP P49189
F	-3	SER	-	expression tag	UNP P49189
F	-2	GLN	-	expression tag	UNP P49189
F	-1	ASP	-	expression tag	UNP P49189
F	0	PRO	-	expression tag	UNP P49189
G	-13	MET	-	initiating methionine	UNP P49189
G	-12	GLY	-	expression tag	UNP P49189
G	-11	SER	-	expression tag	UNP P49189
G	-10	SER	-	expression tag	UNP P49189
G	-9	HIS	-	expression tag	UNP P49189
G	-8	HIS	-	expression tag	UNP P49189
G	-7	HIS	-	expression tag	UNP P49189
G	-6	HIS	-	expression tag	UNP P49189
G	-5	HIS	-	expression tag	UNP P49189
G	-4	HIS	-	expression tag	UNP P49189
G	-3	SER	-	expression tag	UNP P49189
G	-2	GLN	-	expression tag	UNP P49189
G	-1	ASP	-	expression tag	UNP P49189

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	PRO	-	expression tag	UNP P49189
H	-13	MET	-	initiating methionine	UNP P49189
H	-12	GLY	-	expression tag	UNP P49189
H	-11	SER	-	expression tag	UNP P49189
H	-10	SER	-	expression tag	UNP P49189
H	-9	HIS	-	expression tag	UNP P49189
H	-8	HIS	-	expression tag	UNP P49189
H	-7	HIS	-	expression tag	UNP P49189
H	-6	HIS	-	expression tag	UNP P49189
H	-5	HIS	-	expression tag	UNP P49189
H	-4	HIS	-	expression tag	UNP P49189
H	-3	SER	-	expression tag	UNP P49189
H	-2	GLN	-	expression tag	UNP P49189
H	-1	ASP	-	expression tag	UNP P49189
H	0	PRO	-	expression tag	UNP P49189

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).



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

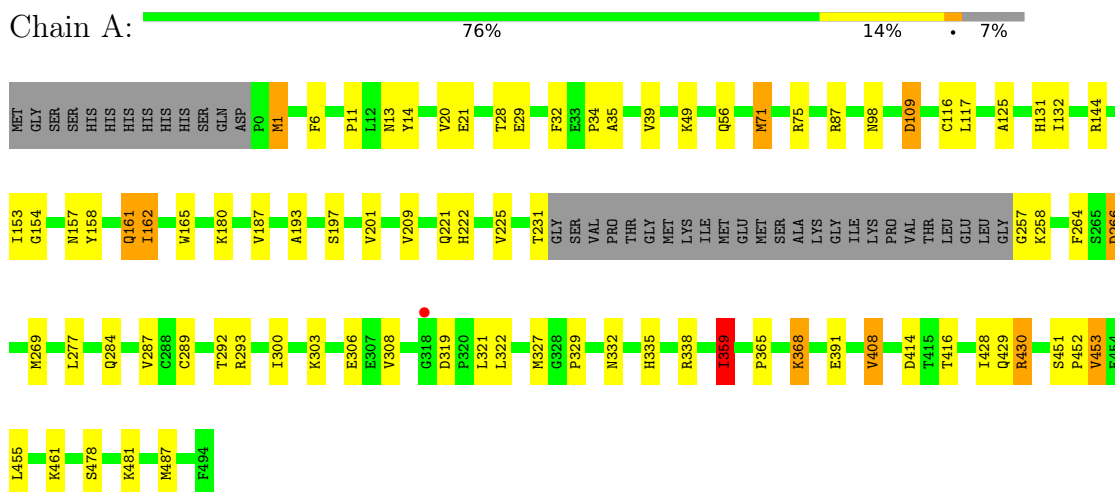
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	O 1	0	0
3	B	4	Total 4	O 4	0	0
3	C	5	Total 5	O 5	0	0
3	D	7	Total 7	O 7	0	0
3	E	3	Total 3	O 3	0	0
3	F	4	Total 4	O 4	0	0
3	H	2	Total 2	O 2	0	0

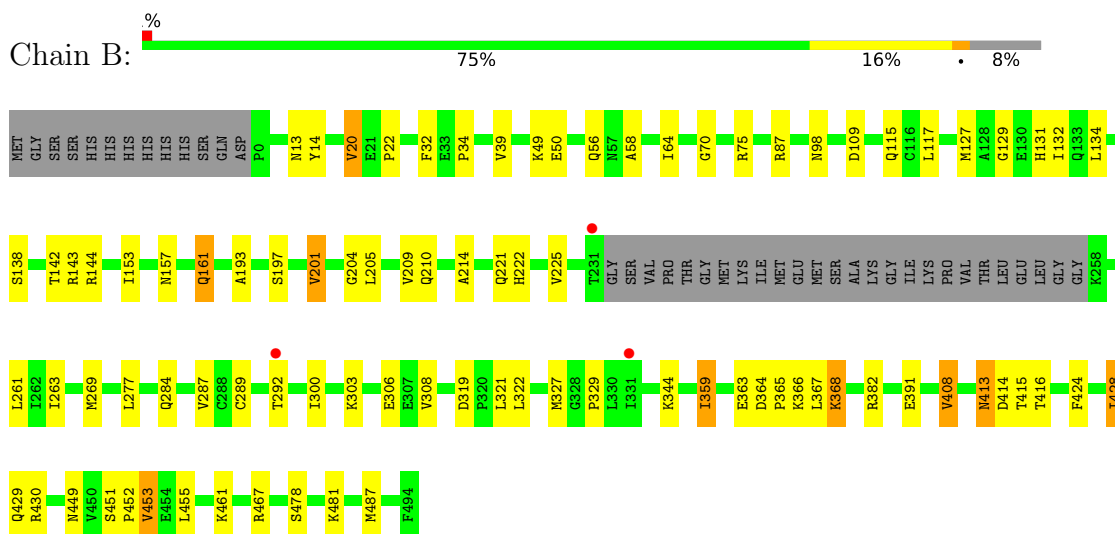
3 Residue-property plots

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

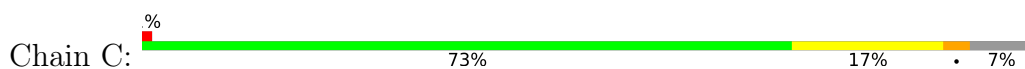
- Molecule 1: 4-trimethylaminobutyraldehyde dehydrogenase

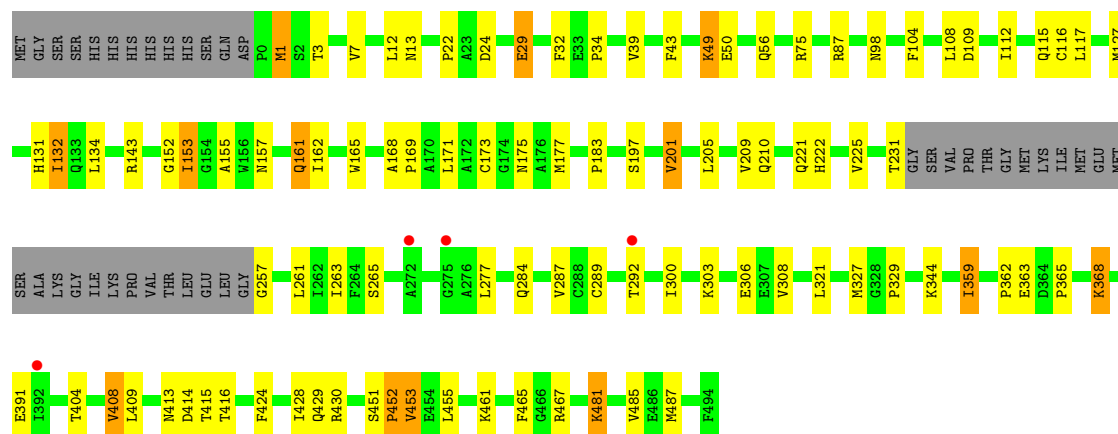


- Molecule 1: 4-trimethylaminobutyraldehyde dehydrogenase



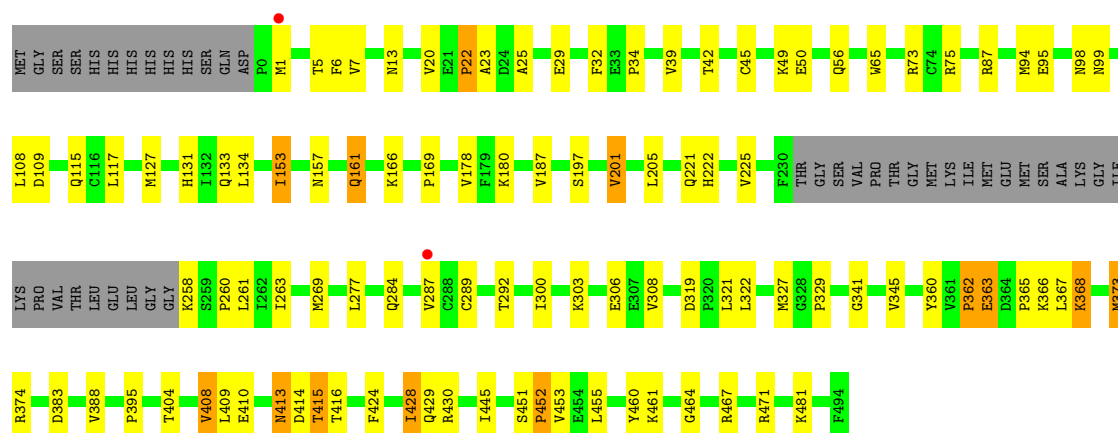
- Molecule 1: 4-trimethylaminobutyraldehyde dehydrogenase





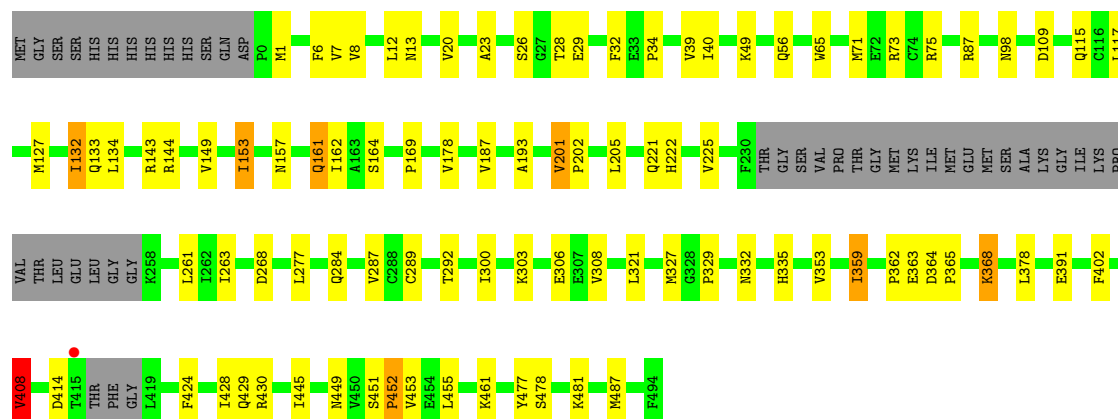
- Molecule 1: 4-trimethylaminobutylaldehyde dehydrogenase

Chain D: 72% 18% 8%



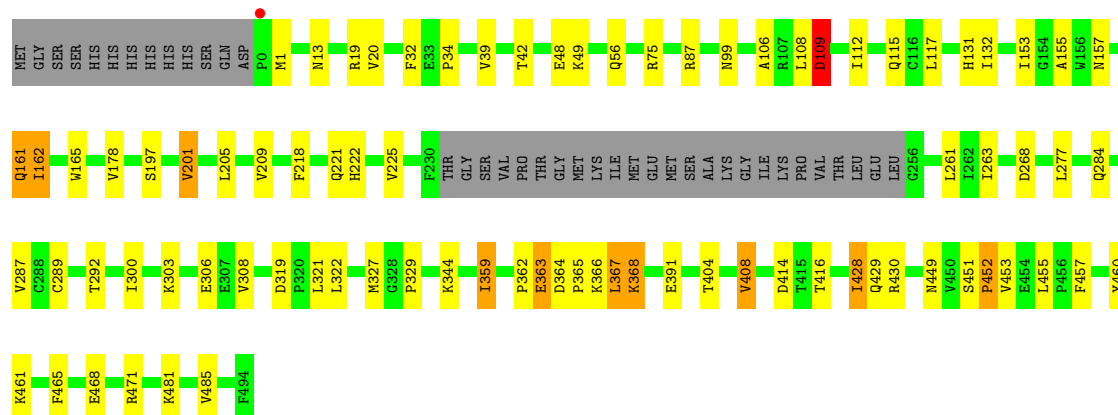
- Molecule 1: 4-trimethylaminobutylaldehyde dehydrogenase

Chain E: 73% 17% 8%



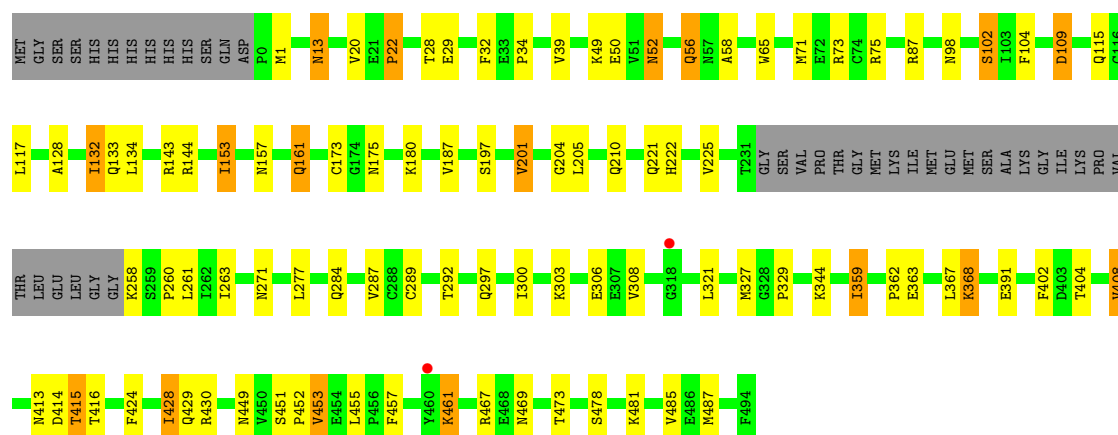
- Molecule 1: 4-trimethylaminobutylaldehyde dehydrogenase

Chain F:  76% 14% 7%



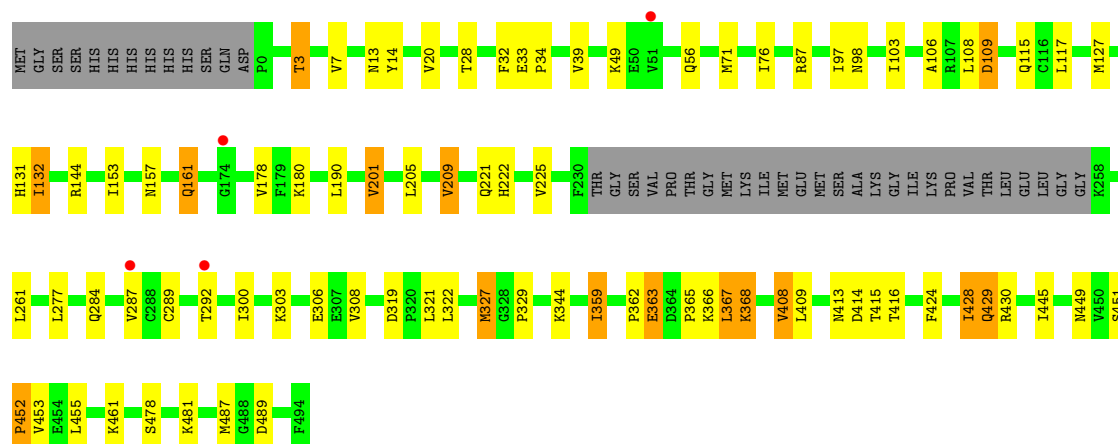
• Molecule 1: 4-trimethylaminobutyraldehyde dehydrogenase

Chain G:  73% 16% 8%



• Molecule 1: 4-trimethylaminobutyraldehyde dehydrogenase

Chain H:  76% 13% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	113.73Å 167.45Å 116.14Å 90.00° 90.94° 90.00°	Depositor
Resolution (Å)	49.71 – 2.89 49.71 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.71-2.89) 99.5 (49.71-2.89)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.91Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.179 , 0.213 0.193 , 0.227	Depositor DCC
R_{free} test set	4819 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	79.7	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 84.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.016 for h,-k,-l 0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	28772	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.90	1/3672 (0.0%)	1.41	14/4966 (0.3%)
1	B	0.86	0/3668	1.39	18/4961 (0.4%)
1	C	0.86	2/3672 (0.1%)	1.42	22/4966 (0.4%)
1	D	0.90	1/3669 (0.0%)	1.42	24/4961 (0.5%)
1	E	0.87	1/3637 (0.0%)	1.42	20/4917 (0.4%)
1	F	0.87	1/3669 (0.0%)	1.41	20/4961 (0.4%)
1	G	0.88	2/3668 (0.1%)	1.41	18/4961 (0.4%)
1	H	0.86	1/3661 (0.0%)	1.39	17/4951 (0.3%)
All	All	0.88	9/29316 (0.0%)	1.41	153/39644 (0.4%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	127	MET	SD-CE	5.95	1.94	1.79
1	E	153	ILE	CG1-CD1	-5.93	1.28	1.51
1	D	153	ILE	CG1-CD1	-5.71	1.29	1.51
1	F	1	MET	SD-CE	5.40	1.93	1.79
1	H	127	MET	SD-CE	5.39	1.93	1.79
1	G	71	MET	SD-CE	5.24	1.92	1.79
1	C	153	ILE	CG1-CD1	-5.12	1.31	1.51
1	A	359	ILE	CG1-CD1	5.05	1.71	1.51
1	G	153	ILE	CG1-CD1	-5.02	1.32	1.51

All (153) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	ASP	CA-CB-CG	8.84	121.44	112.60
1	E	362	PRO	CA-C-N	8.42	131.92	120.38
1	E	362	PRO	C-N-CA	8.42	131.92	120.38
1	G	362	PRO	CA-C-N	8.05	131.41	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	362	PRO	C-N-CA	8.05	131.41	120.38
1	E	187	VAL	N-CA-C	7.96	117.90	110.42
1	H	362	PRO	CA-C-N	7.80	131.06	120.38
1	H	362	PRO	C-N-CA	7.80	131.06	120.38
1	A	187	VAL	N-CA-C	7.79	117.75	110.42
1	D	98	ASN	CA-CB-CG	7.68	120.28	112.60
1	G	187	VAL	N-CA-C	7.53	117.50	110.42
1	G	98	ASN	CA-CB-CG	7.47	120.08	112.60
1	A	201	VAL	N-CA-CB	7.36	117.57	109.99
1	G	413	ASN	CA-CB-CG	7.34	119.94	112.60
1	F	109	ASP	CA-CB-CG	7.29	119.89	112.60
1	A	98	ASN	CA-CB-CG	7.03	119.63	112.60
1	C	362	PRO	CA-C-N	6.99	130.59	120.38
1	C	362	PRO	C-N-CA	6.99	130.59	120.38
1	E	115	GLN	CA-C-N	6.98	129.52	120.44
1	E	115	GLN	C-N-CA	6.98	129.52	120.44
1	E	98	ASN	CA-CB-CG	6.93	119.53	112.60
1	C	98	ASN	CA-CB-CG	6.91	119.51	112.60
1	F	362	PRO	CA-C-N	6.91	130.46	120.38
1	F	362	PRO	C-N-CA	6.91	130.46	120.38
1	H	98	ASN	CA-CB-CG	6.75	119.35	112.60
1	C	115	GLN	CA-C-N	6.72	129.17	120.44
1	C	115	GLN	C-N-CA	6.72	129.17	120.44
1	B	98	ASN	CA-CB-CG	6.64	119.24	112.60
1	G	115	GLN	CA-C-N	6.64	129.07	120.44
1	G	115	GLN	C-N-CA	6.64	129.07	120.44
1	H	115	GLN	CA-C-N	6.60	129.02	120.44
1	H	115	GLN	C-N-CA	6.60	129.02	120.44
1	H	109	ASP	CA-CB-CG	6.54	119.14	112.60
1	A	257	GLY	CA-C-N	6.52	133.22	122.07
1	A	257	GLY	C-N-CA	6.52	133.22	122.07
1	D	187	VAL	N-CA-C	6.40	116.56	110.74
1	D	408	VAL	N-CA-CB	6.33	117.54	110.51
1	H	209	VAL	N-CA-CB	6.32	119.84	111.25
1	E	26	SER	N-CA-C	-6.29	104.87	113.30
1	G	109	ASP	CA-CB-CG	6.24	118.84	112.60
1	B	209	VAL	N-CA-CB	6.23	119.75	111.90
1	H	408	VAL	N-CA-CB	6.22	117.42	110.51
1	D	115	GLN	CA-C-N	6.21	128.51	120.44
1	D	115	GLN	C-N-CA	6.21	128.51	120.44
1	C	257	GLY	CA-C-N	6.21	132.69	122.07
1	C	257	GLY	C-N-CA	6.21	132.69	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	362	PRO	CA-C-N	6.18	133.35	121.54
1	D	362	PRO	C-N-CA	6.18	133.35	121.54
1	B	115	GLN	CA-C-N	6.17	128.47	120.44
1	B	115	GLN	C-N-CA	6.17	128.47	120.44
1	G	408	VAL	N-CA-CB	6.13	117.31	110.51
1	E	109	ASP	CA-CB-CG	6.09	118.69	112.60
1	A	408	VAL	N-CA-CB	6.08	117.25	110.51
1	C	408	VAL	N-CA-CB	6.03	117.21	110.51
1	B	408	VAL	N-CA-CB	5.90	117.06	110.51
1	A	391	GLU	CB-CG-CD	5.88	122.60	112.60
1	F	209	VAL	N-CA-CB	5.88	118.87	111.46
1	A	209	VAL	N-CA-CB	5.88	119.52	111.41
1	E	162	ILE	CA-C-N	5.87	128.07	120.44
1	E	162	ILE	C-N-CA	5.87	128.07	120.44
1	E	408	VAL	N-CA-CB	5.86	116.77	110.62
1	C	452	PRO	CA-C-N	5.85	128.14	120.60
1	C	452	PRO	C-N-CA	5.85	128.14	120.60
1	F	408	VAL	N-CA-CB	5.78	116.93	110.51
1	F	115	GLN	CA-C-N	5.78	127.95	120.44
1	F	115	GLN	C-N-CA	5.78	127.95	120.44
1	D	452	PRO	CA-C-N	5.76	128.03	120.60
1	D	452	PRO	C-N-CA	5.76	128.03	120.60
1	E	193	ALA	CA-C-N	5.75	127.99	120.28
1	E	193	ALA	C-N-CA	5.75	127.99	120.28
1	D	413	ASN	CA-C-N	5.74	132.51	121.54
1	D	413	ASN	C-N-CA	5.74	132.51	121.54
1	E	178	VAL	N-CA-C	-5.73	99.38	107.75
1	G	22	PRO	CA-C-N	5.70	127.92	120.28
1	G	22	PRO	C-N-CA	5.70	127.92	120.28
1	C	197	SER	CA-C-N	5.67	127.88	120.28
1	C	197	SER	C-N-CA	5.67	127.88	120.28
1	D	108	LEU	CA-C-N	5.66	127.79	120.44
1	D	108	LEU	C-N-CA	5.66	127.79	120.44
1	B	64	ILE	CA-C-N	5.58	128.02	120.38
1	B	64	ILE	C-N-CA	5.58	128.02	120.38
1	B	453	VAL	CA-C-N	5.47	128.37	120.38
1	B	453	VAL	C-N-CA	5.47	128.37	120.38
1	D	22	PRO	CA-C-N	5.46	127.59	120.28
1	D	22	PRO	C-N-CA	5.46	127.59	120.28
1	B	214	ALA	CA-C-N	5.43	127.56	120.28
1	B	214	ALA	C-N-CA	5.43	127.56	120.28
1	C	453	VAL	CA-C-N	5.42	127.80	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	453	VAL	C-N-CA	5.42	127.80	120.38
1	B	391	GLU	CB-CG-CD	5.42	121.81	112.60
1	B	197	SER	CA-C-N	5.41	127.53	120.28
1	B	197	SER	C-N-CA	5.41	127.53	120.28
1	B	413	ASN	CA-CB-CG	5.40	118.00	112.60
1	B	109	ASP	CA-CB-CG	5.40	118.00	112.60
1	C	209	VAL	N-CA-CB	5.38	118.63	111.64
1	F	452	PRO	CA-C-N	5.37	127.53	120.60
1	F	452	PRO	C-N-CA	5.37	127.53	120.60
1	H	3	THR	CB-CA-C	5.37	118.83	109.65
1	F	268	ASP	CA-C-N	5.37	128.87	120.82
1	F	268	ASP	C-N-CA	5.37	128.87	120.82
1	D	363	GLU	CB-CG-CD	5.32	121.64	112.60
1	D	197	SER	CA-C-N	5.28	127.35	120.28
1	D	197	SER	C-N-CA	5.28	127.35	120.28
1	H	452	PRO	CA-C-N	5.28	127.41	120.60
1	H	452	PRO	C-N-CA	5.28	127.41	120.60
1	B	70	GLY	CA-C-N	5.27	127.61	120.44
1	B	70	GLY	C-N-CA	5.27	127.61	120.44
1	E	364	ASP	CA-CB-CG	5.25	117.84	112.60
1	F	391	GLU	CB-CG-CD	5.23	121.49	112.60
1	D	410	GLU	CA-C-N	5.22	128.24	120.31
1	D	410	GLU	C-N-CA	5.22	128.24	120.31
1	G	404	THR	CA-C-N	5.20	127.25	120.28
1	G	404	THR	C-N-CA	5.20	127.25	120.28
1	E	268	ASP	CA-C-N	5.20	128.62	120.82
1	E	268	ASP	C-N-CA	5.20	128.62	120.82
1	F	108	LEU	CA-C-N	5.20	127.20	120.44
1	F	108	LEU	C-N-CA	5.20	127.20	120.44
1	H	178	VAL	N-CA-C	-5.20	100.27	107.80
1	H	108	LEU	CA-C-N	5.19	127.19	120.44
1	H	108	LEU	C-N-CA	5.19	127.19	120.44
1	A	109	ASP	CA-CB-CG	5.19	117.79	112.60
1	E	391	GLU	CB-CG-CD	5.17	121.40	112.60
1	F	42	THR	N-CA-C	-5.15	99.01	108.02
1	G	197	SER	CA-C-N	5.15	127.18	120.28
1	G	197	SER	C-N-CA	5.15	127.18	120.28
1	G	391	GLU	CB-CG-CD	5.13	121.33	112.60
1	A	197	SER	CA-C-N	5.12	127.14	120.28
1	A	197	SER	C-N-CA	5.12	127.14	120.28
1	C	391	GLU	CB-CG-CD	5.11	121.28	112.60
1	F	178	VAL	N-CA-C	-5.11	100.39	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	453	VAL	CA-C-N	5.10	127.82	120.38
1	G	453	VAL	C-N-CA	5.10	127.82	120.38
1	E	452	PRO	CA-C-N	5.07	127.68	120.53
1	E	452	PRO	C-N-CA	5.07	127.68	120.53
1	F	404	THR	CA-C-N	5.07	127.07	120.28
1	F	404	THR	C-N-CA	5.07	127.07	120.28
1	C	108	LEU	CA-C-N	5.06	127.02	120.44
1	C	108	LEU	C-N-CA	5.06	127.02	120.44
1	D	178	VAL	N-CA-C	-5.06	100.46	107.80
1	C	404	THR	CA-C-N	5.06	127.06	120.28
1	C	404	THR	C-N-CA	5.06	127.06	120.28
1	C	104	PHE	CA-C-N	5.05	127.47	120.29
1	C	104	PHE	C-N-CA	5.05	127.47	120.29
1	A	453	VAL	CA-C-N	5.03	127.27	120.38
1	A	453	VAL	C-N-CA	5.03	127.27	120.38
1	H	109	ASP	CB-CA-C	-5.03	102.99	110.88
1	D	404	THR	CA-C-N	5.02	127.01	120.28
1	D	404	THR	C-N-CA	5.02	127.01	120.28
1	D	42	THR	N-CA-C	-5.02	99.23	108.02
1	F	197	SER	CA-C-N	5.02	127.01	120.28
1	F	197	SER	C-N-CA	5.02	127.01	120.28
1	H	103	ILE	CA-C-N	5.01	127.27	120.65
1	H	103	ILE	C-N-CA	5.01	127.27	120.65

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	3600	0	3574	31	0
1	B	3596	0	3571	31	0
1	C	3600	0	3574	37	0
1	D	3594	0	3573	40	0
1	E	3567	0	3544	36	0
1	F	3597	0	3570	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	3596	0	3571	33	0
1	H	3589	0	3564	32	0
2	B	7	0	10	0	0
3	A	1	0	0	0	0
3	B	4	0	0	0	0
3	C	5	0	0	0	0
3	D	7	0	0	0	0
3	E	3	0	0	0	0
3	F	4	0	0	0	0
3	H	2	0	0	0	0
All	All	28772	0	28551	243	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 4.

All (243) 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:154:GLY:HA2	1:A:162:ILE:HD11	1.36	1.05
1:E:132:ILE:HD13	1:F:457:PHE:CZ	2.12	0.84
1:A:258:LYS:HD2	1:A:293:ARG:HG3	1.61	0.81
1:C:152:GLY:HA3	1:C:177:MET:HE3	1.61	0.80
1:E:201:VAL:HG13	1:E:205:LEU:HB3	1.62	0.79
1:A:71:MET:HE2	1:A:125:ALA:HB1	1.65	0.79
1:G:457:PHE:CZ	1:H:132:ILE:HD13	2.20	0.77
1:D:1[B]:MET:HE2	1:D:6:PHE:HB2	1.69	0.73
1:F:201:VAL:HG13	1:F:205:LEU:HB3	1.75	0.69
1:D:260:PRO:HD3	1:D:415:THR:HG21	1.75	0.68
1:G:260:PRO:HD3	1:G:415:THR:HG21	1.74	0.68
1:B:428:ILE:HD13	1:C:487:MET:HE1	1.74	0.68
1:B:201:VAL:HG13	1:B:205:LEU:HB3	1.77	0.66
1:G:428:ILE:HD13	1:G:428:ILE:H	1.61	0.66
1:C:201:VAL:HG13	1:C:205:LEU:HB3	1.76	0.66
1:E:487:MET:HE1	1:H:428:ILE:HD13	1.78	0.65
1:F:300:ILE:HA	1:F:303:LYS:HE2	1.79	0.65
1:B:300:ILE:HA	1:B:303:LYS:HE2	1.80	0.64
1:G:300:ILE:HA	1:G:303:LYS:HE2	1.80	0.64
1:D:201:VAL:HG13	1:D:205:LEU:HB3	1.78	0.64
1:A:300:ILE:HA	1:A:303:LYS:HE2	1.80	0.63
1:C:300:ILE:HA	1:C:303:LYS:HE2	1.80	0.63
1:G:102:SER:HB2	1:G:104:PHE:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:300:ILE:HA	1:E:303:LYS:HE2	1.80	0.62
1:D:300:ILE:HA	1:D:303:LYS:HE2	1.81	0.62
1:F:161:GLN:OE1	1:F:452:PRO:HG2	1.99	0.62
1:C:29:GLU:HG2	1:C:183:PRO:HB2	1.81	0.62
1:B:143:ARG:HD3	1:D:133:GLN:HG3	1.82	0.62
1:A:487:MET:HE1	1:D:428:ILE:HD13	1.81	0.61
1:H:300:ILE:HA	1:H:303:LYS:HE2	1.82	0.61
1:C:143:ARG:HH21	1:C:481:LYS:HD2	1.65	0.61
1:B:13:ASN:HB2	1:B:20:VAL:O	2.02	0.59
1:A:332:ASN:OD1	1:A:335:HIS:HB2	2.02	0.59
1:G:201:VAL:HG13	1:G:205:LEU:HB3	1.84	0.59
1:G:32:PHE:CE1	1:G:39:VAL:HG22	2.37	0.59
1:F:131:HIS:HB2	1:H:131:HIS:HB2	1.86	0.58
1:G:58:ALA:HA	1:G:204:GLY:O	2.04	0.57
1:D:34:PRO:HB2	1:D:329:PRO:HG2	1.87	0.56
1:D:362:PRO:HG2	1:D:367:LEU:HB3	1.86	0.56
1:A:14:TYR:CD2	1:A:193:ALA:HB1	2.41	0.56
1:F:366:LYS:HG3	1:F:367:LEU:HD13	1.87	0.56
1:F:165:TRP:HH2	1:F:465:PHE:HE2	1.53	0.56
1:E:8:VAL:HG11	1:E:12:LEU:HD21	1.89	0.55
1:H:201:VAL:HG13	1:H:205:LEU:HB3	1.88	0.55
1:E:34:PRO:HB2	1:E:329:PRO:HG2	1.88	0.55
1:A:264:PHE:HB3	1:A:430:ARG:HH21	1.71	0.55
1:A:34:PRO:HB2	1:A:329:PRO:HG2	1.89	0.55
1:B:131:HIS:HB2	1:D:131:HIS:HB2	1.88	0.54
1:A:359:ILE:H	1:A:359:ILE:HD12	1.72	0.54
1:G:297:GLN:HG3	1:G:402:PHE:CE1	2.43	0.54
1:B:364:ASP:HB3	1:B:367:LEU:HD22	1.90	0.54
1:D:161:GLN:OE1	1:D:452:PRO:HG2	2.08	0.54
1:F:365:PRO:O	1:F:368:LYS:HB2	2.08	0.54
1:D:1[B]:MET:HE1	1:D:5:THR:OG1	2.07	0.54
1:H:409:LEU:O	1:H:413:ASN:HB2	2.08	0.53
1:H:34:PRO:HB2	1:H:329:PRO:HG2	1.89	0.53
1:B:451:SER:HB2	1:B:455:LEU:HD12	1.91	0.53
1:E:359:ILE:HD12	1:E:359:ILE:H	1.73	0.53
1:F:165:TRP:CH2	1:F:465:PHE:HE2	2.26	0.53
1:G:132:ILE:HG22	1:G:134:LEU:CD1	2.39	0.53
1:F:34:PRO:HB2	1:F:329:PRO:HG2	1.90	0.53
1:G:359:ILE:HD12	1:G:359:ILE:H	1.74	0.53
1:A:365:PRO:O	1:A:368:LYS:HB2	2.10	0.52
1:E:161:GLN:OE1	1:E:452:PRO:HG2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:GLN:HA	1:B:327:MET:HE2	1.90	0.52
1:E:132:ILE:HD11	1:G:128:ALA:HB2	1.90	0.52
1:C:284:GLN:HA	1:C:327:MET:HE2	1.91	0.52
1:E:451:SER:HB2	1:E:455:LEU:HD12	1.92	0.52
1:E:132:ILE:HG22	1:E:134:LEU:CD1	2.40	0.52
1:E:428:ILE:HD13	1:H:487:MET:HE1	1.92	0.52
1:A:158:TYR:HB2	1:A:162:ILE:HG23	1.92	0.51
1:H:365:PRO:O	1:H:368:LYS:HB2	2.11	0.51
1:A:284:GLN:HA	1:A:327:MET:HE2	1.92	0.51
1:C:132:ILE:HG22	1:C:134:LEU:CD1	2.39	0.51
1:C:451:SER:HB2	1:C:455:LEU:HD12	1.92	0.51
1:F:284:GLN:HA	1:F:327:MET:HE2	1.91	0.51
1:H:106:ALA:O	1:H:109:ASP:HB2	2.11	0.51
1:B:34:PRO:HB2	1:B:329:PRO:HG2	1.92	0.51
1:H:161:GLN:OE1	1:H:452:PRO:HG2	2.11	0.51
1:B:161:GLN:OE1	1:B:452:PRO:HG2	2.11	0.51
1:C:34:PRO:HB2	1:C:329:PRO:HG2	1.92	0.51
1:D:366:LYS:HG3	1:D:367:LEU:HD13	1.92	0.51
1:E:284:GLN:HA	1:E:327:MET:HE2	1.93	0.51
1:C:161:GLN:OE1	1:C:452:PRO:HG2	2.11	0.51
1:G:34:PRO:HB2	1:G:329:PRO:HG2	1.92	0.51
1:C:162:ILE:HD11	1:C:231:THR:OG1	2.10	0.51
1:F:165:TRP:HH2	1:F:465:PHE:CE2	2.30	0.50
1:F:451:SER:HB2	1:F:455:LEU:HD12	1.93	0.50
1:B:359:ILE:HD12	1:B:359:ILE:H	1.76	0.50
1:G:284:GLN:HA	1:G:327:MET:HE2	1.92	0.50
1:B:428:ILE:HD12	1:D:428:ILE:HD12	1.93	0.50
1:H:284:GLN:HA	1:H:327:MET:HE2	1.93	0.50
1:F:359:ILE:H	1:F:359:ILE:HD12	1.75	0.50
1:D:284:GLN:HA	1:D:327:MET:HE2	1.92	0.50
1:E:332:ASN:OD1	1:E:335:HIS:HB2	2.12	0.50
1:C:152:GLY:HA3	1:C:177:MET:CE	2.38	0.50
1:H:451:SER:HB2	1:H:455:LEU:HD12	1.94	0.49
1:B:366:LYS:HG3	1:B:367:LEU:HD13	1.94	0.49
1:D:166:LYS:O	1:D:169:PRO:HD2	2.13	0.49
1:F:428:ILE:HD13	1:G:487:MET:HE1	1.94	0.49
1:C:24:ASP:HB2	1:C:49:LYS:HB2	1.94	0.49
1:C:409:LEU:O	1:C:413:ASN:HB2	2.13	0.49
1:C:365:PRO:O	1:C:368:LYS:HB2	2.12	0.49
1:A:365:PRO:HD2	1:E:23:ALA:HB1	1.95	0.49
1:B:58:ALA:HA	1:B:204:GLY:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:161:GLN:OE1	1:G:452:PRO:HG2	2.12	0.49
1:G:451:SER:HB2	1:G:455:LEU:HD12	1.94	0.49
1:E:201:VAL:HG22	1:E:205:LEU:HD23	1.93	0.49
1:H:359:ILE:HD12	1:H:359:ILE:H	1.76	0.48
1:B:487:MET:HE1	1:C:428:ILE:HD13	1.95	0.48
1:G:22:PRO:HB3	1:G:50:GLU:HG2	1.93	0.48
1:H:366:LYS:HG3	1:H:367:LEU:HD13	1.95	0.48
1:B:365:PRO:O	1:B:368:LYS:HB2	2.14	0.48
1:C:359:ILE:HD12	1:C:359:ILE:H	1.77	0.48
1:E:365:PRO:O	1:E:368:LYS:HB2	2.13	0.48
1:E:143:ARG:HD3	1:G:133:GLN:HG3	1.95	0.48
1:A:131:HIS:HB2	1:C:131:HIS:HB2	1.96	0.48
1:E:222:HIS:HB3	1:E:225:VAL:HG23	1.96	0.48
1:H:413:ASN:C	1:H:415:THR:H	2.22	0.48
1:D:451:SER:HB2	1:D:455:LEU:HD12	1.95	0.47
1:F:13:ASN:HB2	1:F:20:VAL:O	2.14	0.47
1:B:413:ASN:C	1:B:415:THR:H	2.23	0.47
1:D:365:PRO:O	1:D:368:LYS:HB2	2.14	0.47
1:C:116:CYS:HB2	1:C:165:TRP:CZ3	2.49	0.47
1:B:210:GLN:HA	1:B:210:GLN:NE2	2.30	0.47
1:D:373:MET:HE2	1:D:395:PRO:HG2	1.97	0.47
1:D:409:LEU:O	1:D:413:ASN:HB2	2.14	0.47
1:D:360:TYR:HB2	1:D:374:ARG:HG3	1.96	0.47
1:E:65:TRP:CE2	1:E:73:ARG:HD2	2.50	0.47
1:A:32:PHE:CE2	1:A:39:VAL:HG22	2.50	0.47
1:A:1:MET:SD	1:A:6:PHE:HB2	2.56	0.46
1:A:71:MET:HE3	1:D:464:GLY:HA2	1.97	0.46
1:E:13:ASN:HB2	1:E:20:VAL:O	2.15	0.46
1:E:201:VAL:CG1	1:E:205:LEU:HB3	2.41	0.46
1:A:158:TYR:O	1:A:162:ILE:HG23	2.15	0.46
1:D:23:ALA:HB1	1:H:365:PRO:HD2	1.97	0.46
1:E:133:GLN:HG3	1:G:143:ARG:HD3	1.97	0.46
1:H:13:ASN:HB2	1:H:20:VAL:O	2.16	0.46
1:F:364:ASP:HB3	1:F:367:LEU:HD22	1.97	0.46
1:G:210:GLN:OE1	1:G:210:GLN:HA	2.16	0.46
1:B:222:HIS:HB3	1:B:225:VAL:HG23	1.97	0.45
1:C:168:ALA:HB3	1:C:169:PRO:HD3	1.98	0.45
1:G:428:ILE:H	1:G:428:ILE:CD1	2.28	0.45
1:B:14:TYR:CD2	1:B:193:ALA:HB1	2.51	0.45
1:D:65:TRP:CE2	1:D:73:ARG:HD2	2.52	0.45
1:F:32:PHE:CE2	1:F:39:VAL:HG22	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:445:ILE:HD12	1:F:485:VAL:HG22	1.98	0.45
1:D:287:VAL:HG12	1:D:289:CYS:H	1.82	0.45
1:F:13:ASN:O	1:F:19:ARG:HA	2.17	0.45
1:C:287:VAL:HG12	1:C:289:CYS:H	1.82	0.45
1:F:201:VAL:CG1	1:F:205:LEU:HB3	2.46	0.45
1:C:413:ASN:C	1:C:415:THR:H	2.25	0.45
1:C:261:LEU:HD11	1:C:424:PHE:HE1	1.81	0.45
1:D:95:GLU:O	1:D:99:ASN:HB3	2.17	0.45
1:F:222:HIS:HB3	1:F:225:VAL:HG23	1.98	0.45
1:G:287:VAL:HG12	1:G:289:CYS:H	1.82	0.45
1:A:451:SER:HB2	1:A:455:LEU:HD22	1.98	0.45
1:C:222:HIS:HB3	1:C:225:VAL:HG23	1.99	0.45
1:E:287:VAL:HG12	1:E:289:CYS:H	1.81	0.45
1:E:487:MET:HE1	1:H:428:ILE:CD1	2.45	0.45
1:C:485:VAL:HG22	1:D:445:ILE:HD12	1.99	0.45
1:G:261:LEU:HD11	1:G:424:PHE:HE1	1.82	0.45
1:E:487:MET:HE2	1:H:429:GLN:HA	1.99	0.44
1:C:210:GLN:HA	1:C:210:GLN:NE2	2.32	0.44
1:F:468:GLU:HG3	1:F:471:ARG:HH22	1.83	0.44
1:B:287:VAL:HG12	1:B:289:CYS:H	1.82	0.44
1:D:1[A]:MET:SD	1:D:94:MET:HE2	2.58	0.44
1:G:222:HIS:HB3	1:G:225:VAL:HG23	1.98	0.44
1:D:261:LEU:HD11	1:D:424:PHE:HE1	1.83	0.44
1:G:144:ARG:HD3	1:G:478:SER:OG	2.18	0.44
1:B:261:LEU:HD11	1:B:424:PHE:HE1	1.83	0.44
1:G:13:ASN:HB2	1:G:20:VAL:O	2.18	0.44
1:G:485:VAL:HG22	1:H:445:ILE:HD12	2.00	0.44
1:H:222:HIS:HB3	1:H:225:VAL:HG23	2.00	0.44
1:B:22:PRO:HB3	1:B:50:GLU:HG2	2.00	0.43
1:A:158:TYR:HB3	1:A:161:GLN:HB3	2.00	0.43
1:G:52:ASN:O	1:G:56:GLN:HB2	2.19	0.43
1:A:116:CYS:HB2	1:A:165:TRP:CZ3	2.53	0.43
1:E:6:PHE:O	1:E:40:ILE:HG22	2.18	0.43
1:A:161:GLN:OE1	1:A:452:PRO:HG2	2.18	0.43
1:E:144:ARG:HD3	1:E:478:SER:OG	2.19	0.43
1:A:287:VAL:HG12	1:A:289:CYS:H	1.82	0.43
1:B:134:LEU:HB2	1:B:138:SER:O	2.19	0.43
1:D:32:PHE:CE2	1:D:39:VAL:HG22	2.53	0.43
1:H:32:PHE:CE2	1:H:39:VAL:HG22	2.53	0.43
1:E:32:PHE:CE2	1:E:39:VAL:HG22	2.53	0.43
1:A:13:ASN:HB2	1:A:20:VAL:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:PHE:CE2	1:B:39:VAL:HG22	2.54	0.43
1:C:32:PHE:CE2	1:C:39:VAL:HG22	2.53	0.43
1:C:173:CYS:HB2	1:C:175:ASN:ND2	2.33	0.43
1:F:287:VAL:HG12	1:F:289:CYS:H	1.83	0.43
1:B:144:ARG:HD3	1:B:478:SER:OG	2.19	0.42
1:D:383:ASP:HA	1:D:388:VAL:HG11	2.01	0.42
1:F:106:ALA:O	1:F:109:ASP:HB2	2.18	0.42
1:D:22:PRO:HB3	1:D:50:GLU:HG2	2.01	0.42
1:H:287:VAL:HG12	1:H:289:CYS:H	1.84	0.42
1:A:269:MET:HE1	1:A:300:ILE:HB	2.02	0.42
1:C:1:MET:H	1:C:1:MET:HG3	1.74	0.42
1:G:173:CYS:HB2	1:G:175:ASN:ND2	2.34	0.42
1:G:469:ASN:O	1:G:473:THR:HG23	2.19	0.42
1:H:261:LEU:HD11	1:H:424:PHE:HE1	1.83	0.42
1:D:25:ALA:HA	1:D:45:CYS:O	2.19	0.42
1:E:353:VAL:HG22	1:E:378:LEU:HD21	2.01	0.42
1:A:222:HIS:HB3	1:A:225:VAL:HG23	2.02	0.42
1:B:364:ASP:HA	1:B:365:PRO:HD3	1.97	0.42
1:D:319:ASP:HB3	1:D:322:LEU:HD22	2.02	0.42
1:D:269:MET:HE1	1:D:300:ILE:HB	2.01	0.42
1:C:134:LEU:HD23	1:D:460:TYR:CD2	2.54	0.41
1:E:202:PRO:HD2	1:E:205:LEU:HD22	2.02	0.41
1:E:402:PHE:CG	1:E:408:VAL:HG13	2.55	0.41
1:A:144:ARG:HD3	1:A:478:SER:OG	2.20	0.41
1:E:261:LEU:HD11	1:E:424:PHE:HE1	1.85	0.41
1:H:33:GLU:HA	1:H:97:ILE:O	2.19	0.41
1:B:129:GLY:HA3	1:B:142:THR:O	2.20	0.41
1:D:222:HIS:HB3	1:D:225:VAL:HG23	2.01	0.41
1:F:155:ALA:H	1:F:162:ILE:HG21	1.85	0.41
1:G:271:ASN:OD1	1:H:489:ASP:HA	2.21	0.41
1:C:155:ALA:H	1:C:162:ILE:HG21	1.86	0.41
1:F:363:GLU:CD	1:F:363:GLU:H	2.28	0.41
1:G:65:TRP:CE2	1:G:73:ARG:HD2	2.55	0.41
1:H:180:LYS:HA	1:H:209:VAL:O	2.21	0.41
1:H:319:ASP:HB3	1:H:322:LEU:HD22	2.01	0.41
1:D:341:GLY:O	1:D:345:VAL:HG23	2.20	0.41
1:A:11:PRO:HB3	1:A:21:GLU:HG2	2.02	0.41
1:D:467:ARG:HG2	1:D:471:ARG:HH21	1.85	0.41
1:A:35:ALA:HA	1:A:329:PRO:HG3	2.03	0.41
1:A:319:ASP:HB3	1:A:322:LEU:HD22	2.03	0.41
1:B:319:ASP:HB3	1:B:322:LEU:HD22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:LEU:HD12	1:C:43:PHE:HB3	2.03	0.41
1:E:169:PRO:HB2	1:E:477:TYR:CD2	2.56	0.41
1:H:14:TYR:HB2	1:H:190:LEU:HD23	2.03	0.41
1:H:363:GLU:H	1:H:363:GLU:CD	2.29	0.41
1:C:22:PRO:HB3	1:C:50:GLU:HG2	2.03	0.41
1:E:134:LEU:HD23	1:F:460:TYR:CD2	2.56	0.41
1:F:319:ASP:HB3	1:F:322:LEU:HD22	2.03	0.41
1:C:165:TRP:HH2	1:C:465:PHE:CE1	2.39	0.40
1:D:1[B]:MET:CE	1:D:6:PHE:HB2	2.47	0.40
1:B:269:MET:HE1	1:B:300:ILE:HB	2.03	0.40
1:C:132:ILE:HG22	1:C:134:LEU:HD11	2.03	0.40
1:F:48:GLU:HB3	1:F:218:PHE:CD1	2.57	0.40
1:C:265:SER:HA	1:C:300:ILE:HG12	2.03	0.40
1:D:13:ASN:HB2	1:D:20:VAL:O	2.20	0.40
1:H:144:ARG:HD3	1:H:478:SER:OG	2.21	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	466/508 (92%)	452 (97%)	13 (3%)	1 (0%)	43	72
1	B	465/508 (92%)	452 (97%)	11 (2%)	2 (0%)	30	59
1	C	466/508 (92%)	452 (97%)	13 (3%)	1 (0%)	43	72
1	D	465/508 (92%)	452 (97%)	11 (2%)	2 (0%)	30	59
1	E	459/508 (90%)	445 (97%)	12 (3%)	2 (0%)	30	59
1	F	466/508 (92%)	454 (97%)	10 (2%)	2 (0%)	30	59
1	G	465/508 (92%)	451 (97%)	11 (2%)	3 (1%)	21	51
1	H	464/508 (91%)	452 (97%)	10 (2%)	2 (0%)	30	59

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	3716/4064 (91%)	3610 (97%)	91 (2%)	15 (0%)	30 59

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	461	LYS
1	F	461	LYS
1	G	368	LYS
1	A	461	LYS
1	D	414	ASP
1	G	414	ASP
1	G	461	LYS
1	H	461	LYS
1	B	461	LYS
1	D	461	LYS
1	F	414	ASP
1	C	414	ASP
1	E	414	ASP
1	B	414	ASP
1	H	414	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	384/416 (92%)	349 (91%)	35 (9%)	9 28
1	B	384/416 (92%)	351 (91%)	33 (9%)	10 30
1	C	384/416 (92%)	347 (90%)	37 (10%)	8 26
1	D	384/416 (92%)	350 (91%)	34 (9%)	9 28
1	E	381/416 (92%)	347 (91%)	34 (9%)	9 28
1	F	383/416 (92%)	348 (91%)	35 (9%)	9 28
1	G	384/416 (92%)	342 (89%)	42 (11%)	6 20
1	H	383/416 (92%)	349 (91%)	34 (9%)	9 28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	3067/3328 (92%)	2783 (91%)	284 (9%)	8 27

All (284) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	28	THR
1	A	29	GLU
1	A	49	LYS
1	A	56	GLN
1	A	71	MET
1	A	75	ARG
1	A	87	ARG
1	A	109	ASP
1	A	117	LEU
1	A	132	ILE
1	A	153	ILE
1	A	157	ASN
1	A	161	GLN
1	A	162	ILE
1	A	180	LYS
1	A	221	GLN
1	A	231	THR
1	A	266	ASP
1	A	277	LEU
1	A	292	THR
1	A	306	GLU
1	A	308	VAL
1	A	321	LEU
1	A	338	ARG
1	A	359	ILE
1	A	368	LYS
1	A	408	VAL
1	A	414	ASP
1	A	416	THR
1	A	428	ILE
1	A	429	GLN
1	A	430	ARG
1	A	453	VAL
1	A	481	LYS
1	B	20	VAL
1	B	49	LYS

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Mol	Chain	Res	Type
1	B	56	GLN
1	B	75	ARG
1	B	87	ARG
1	B	117	LEU
1	B	127	MET
1	B	132	ILE
1	B	153	ILE
1	B	157	ASN
1	B	161	GLN
1	B	201	VAL
1	B	221	GLN
1	B	263	ILE
1	B	277	LEU
1	B	292	THR
1	B	306	GLU
1	B	308	VAL
1	B	321	LEU
1	B	344	LYS
1	B	359	ILE
1	B	363	GLU
1	B	368	LYS
1	B	382	ARG
1	B	408	VAL
1	B	416	THR
1	B	428	ILE
1	B	429	GLN
1	B	430	ARG
1	B	449	ASN
1	B	453	VAL
1	B	467	ARG
1	B	481	LYS
1	C	1	MET
1	C	3	THR
1	C	7	VAL
1	C	13	ASN
1	C	29	GLU
1	C	49	LYS
1	C	56	GLN
1	C	75	ARG
1	C	87	ARG
1	C	109	ASP
1	C	112	ILE

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Mol	Chain	Res	Type
1	C	117	LEU
1	C	132	ILE
1	C	153	ILE
1	C	157	ASN
1	C	161	GLN
1	C	171	LEU
1	C	201	VAL
1	C	221	GLN
1	C	263	ILE
1	C	277	LEU
1	C	292	THR
1	C	306	GLU
1	C	308	VAL
1	C	321	LEU
1	C	344	LYS
1	C	359	ILE
1	C	363	GLU
1	C	368	LYS
1	C	408	VAL
1	C	416	THR
1	C	429	GLN
1	C	430	ARG
1	C	453	VAL
1	C	461	LYS
1	C	467	ARG
1	C	481	LYS
1	D	7	VAL
1	D	29	GLU
1	D	49	LYS
1	D	56	GLN
1	D	75	ARG
1	D	87	ARG
1	D	109	ASP
1	D	117	LEU
1	D	127	MET
1	D	134	LEU
1	D	153	ILE
1	D	157	ASN
1	D	161	GLN
1	D	180	LYS
1	D	201	VAL
1	D	221	GLN

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Mol	Chain	Res	Type
1	D	258	LYS
1	D	263	ILE
1	D	277	LEU
1	D	292	THR
1	D	306	GLU
1	D	308	VAL
1	D	321	LEU
1	D	363	GLU
1	D	368	LYS
1	D	373	MET
1	D	408	VAL
1	D	415	THR
1	D	416	THR
1	D	428	ILE
1	D	429	GLN
1	D	430	ARG
1	D	453	VAL
1	D	481	LYS
1	E	1	MET
1	E	7	VAL
1	E	28	THR
1	E	29	GLU
1	E	49	LYS
1	E	56	GLN
1	E	71	MET
1	E	75	ARG
1	E	87	ARG
1	E	117	LEU
1	E	127	MET
1	E	132	ILE
1	E	149	VAL
1	E	153	ILE
1	E	157	ASN
1	E	161	GLN
1	E	164	SER
1	E	201	VAL
1	E	221	GLN
1	E	263	ILE
1	E	277	LEU
1	E	292	THR
1	E	306	GLU
1	E	308	VAL

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Mol	Chain	Res	Type
1	E	321	LEU
1	E	359	ILE
1	E	363	GLU
1	E	368	LYS
1	E	408	VAL
1	E	429	GLN
1	E	430	ARG
1	E	449	ASN
1	E	453	VAL
1	E	481	LYS
1	F	49	LYS
1	F	56	GLN
1	F	75	ARG
1	F	87	ARG
1	F	99	ASN
1	F	109	ASP
1	F	112	ILE
1	F	117	LEU
1	F	132	ILE
1	F	153	ILE
1	F	157	ASN
1	F	161	GLN
1	F	162	ILE
1	F	201	VAL
1	F	221	GLN
1	F	261	LEU
1	F	263	ILE
1	F	277	LEU
1	F	292	THR
1	F	306	GLU
1	F	308	VAL
1	F	321	LEU
1	F	344	LYS
1	F	359	ILE
1	F	363	GLU
1	F	367	LEU
1	F	368	LYS
1	F	408	VAL
1	F	416	THR
1	F	428	ILE
1	F	429	GLN
1	F	430	ARG

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Mol	Chain	Res	Type
1	F	449	ASN
1	F	453	VAL
1	F	481	LYS
1	G	1	MET
1	G	13	ASN
1	G	28	THR
1	G	29	GLU
1	G	49	LYS
1	G	52	ASN
1	G	56	GLN
1	G	75	ARG
1	G	87	ARG
1	G	102	SER
1	G	109	ASP
1	G	117	LEU
1	G	132	ILE
1	G	153	ILE
1	G	157	ASN
1	G	161	GLN
1	G	180	LYS
1	G	201	VAL
1	G	221	GLN
1	G	258	LYS
1	G	263	ILE
1	G	277	LEU
1	G	292	THR
1	G	306	GLU
1	G	308	VAL
1	G	321	LEU
1	G	344	LYS
1	G	359	ILE
1	G	363	GLU
1	G	367	LEU
1	G	368	LYS
1	G	408	VAL
1	G	415	THR
1	G	416	THR
1	G	428	ILE
1	G	429	GLN
1	G	430	ARG
1	G	449	ASN
1	G	453	VAL

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Mol	Chain	Res	Type
1	G	461	LYS
1	G	467	ARG
1	G	481	LYS
1	H	3	THR
1	H	7	VAL
1	H	28	THR
1	H	49	LYS
1	H	56	GLN
1	H	71	MET
1	H	76	ILE
1	H	87	ARG
1	H	117	LEU
1	H	132	ILE
1	H	153	ILE
1	H	157	ASN
1	H	161	GLN
1	H	201	VAL
1	H	221	GLN
1	H	277	LEU
1	H	292	THR
1	H	306	GLU
1	H	308	VAL
1	H	321	LEU
1	H	327	MET
1	H	344	LYS
1	H	359	ILE
1	H	363	GLU
1	H	367	LEU
1	H	368	LYS
1	H	408	VAL
1	H	416	THR
1	H	428	ILE
1	H	429	GLN
1	H	430	ARG
1	H	449	ASN
1	H	453	VAL
1	H	481	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN

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Mol	Chain	Res	Type
1	A	157	ASN
1	A	270	ASN
1	A	335	HIS
1	A	429	GLN
1	B	67	GLN
1	B	157	ASN
1	B	270	ASN
1	B	449	ASN
1	C	157	ASN
1	C	270	ASN
1	C	290	ASN
1	D	56	GLN
1	D	157	ASN
1	D	290	ASN
1	D	380	ASN
1	E	131	HIS
1	E	157	ASN
1	E	270	ASN
1	E	429	GLN
1	E	449	ASN
1	F	56	GLN
1	F	210	GLN
1	F	270	ASN
1	F	449	ASN
1	G	56	GLN
1	G	157	ASN
1	G	270	ASN
1	G	429	GLN
1	G	449	ASN
1	H	157	ASN
1	H	270	ASN
1	H	449	ASN
1	H	469	ASN

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

1 ligand is 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	PEG	B	501	-	6,6,6	0.44	0	5,5,5	0.35	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PEG	B	501	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

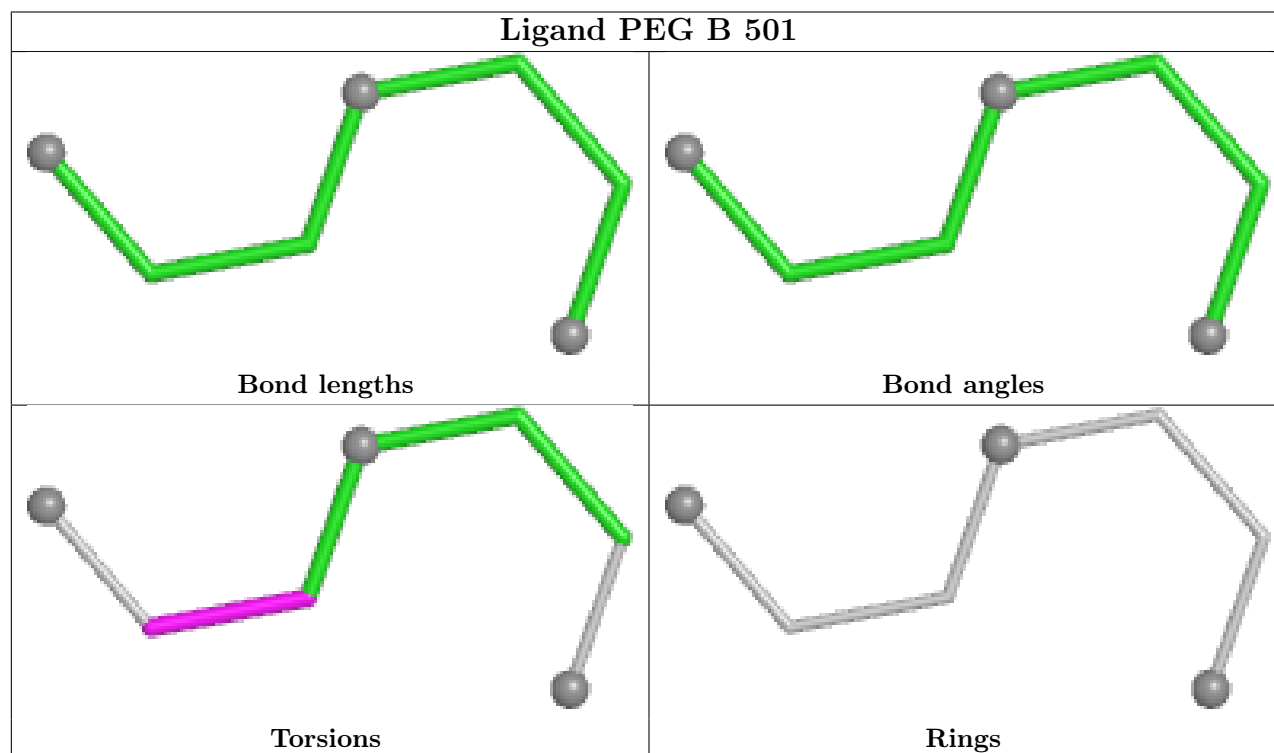
Mol	Chain	Res	Type	Atoms
2	B	501	PEG	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	470/508 (92%)	-0.26	1 (0%) 91 89	54, 78, 112, 127	0
1	B	469/508 (92%)	-0.24	3 (0%) 85 81	53, 87, 116, 132	0
1	C	470/508 (92%)	-0.15	4 (0%) 81 75	60, 99, 135, 157	0
1	D	468/508 (92%)	-0.34	2 (0%) 88 85	53, 75, 107, 143	1 (0%)
1	E	465/508 (91%)	-0.27	1 (0%) 91 89	54, 86, 133, 167	0
1	F	470/508 (92%)	-0.24	1 (0%) 91 89	59, 88, 124, 148	0
1	G	469/508 (92%)	-0.25	2 (0%) 88 85	58, 87, 120, 137	0
1	H	468/508 (92%)	-0.20	4 (0%) 81 75	61, 90, 123, 141	0
All	All	3749/4064 (92%)	-0.24	18 (0%) 87 83	53, 86, 124, 167	1 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	174	GLY	4.8
1	E	415	THR	3.4
1	C	392	ILE	2.9
1	B	231	THR	2.9
1	A	318	GLY	2.5
1	H	51	VAL	2.5
1	H	292	THR	2.5
1	B	331	ILE	2.4
1	C	272	ALA	2.3
1	D	1[A]	MET	2.2
1	F	0	PRO	2.2
1	G	460	TYR	2.1
1	B	292	THR	2.0
1	C	275	GLY	2.0
1	G	318	GLY	2.0
1	C	292	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	287	VAL	2.0
1	H	287	VAL	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](#)

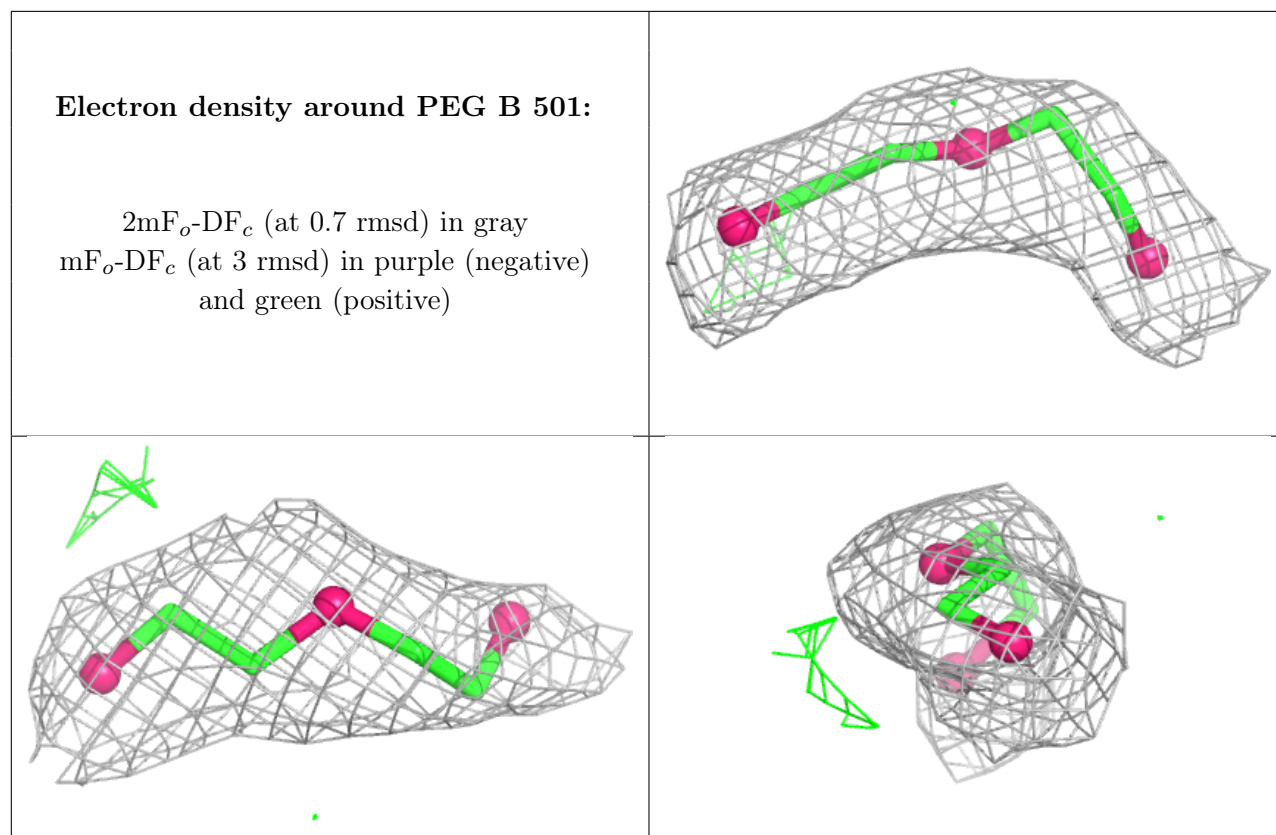
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	PEG	B	501	7/7	0.88	0.10	92,94,97,97	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.