



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:54 PM UTC

PDB ID : 7QNM / pdb_00007qnm
Title : Crystallization and structural analyses of ZgHAD, a L-2-haloacid dehalogenase from the marine Flavobacterium Zobellia galactanivorans
Authors : Grigorian, E.; Roret, T.; Leblanc, C.; Delage, L.; Czjzek, M.
Deposited on : 2021-12-21
Resolution : 2.73 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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
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

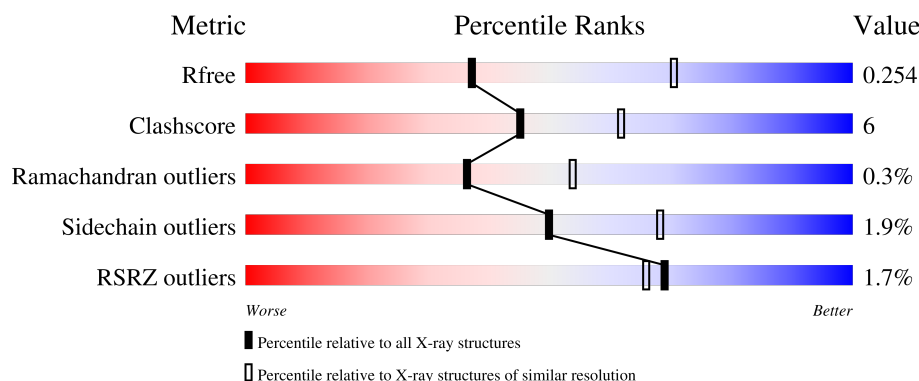
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	227	0% 87% 10% ...
1	B	227	2% 87% 8% ...
1	C	227	0% 80% 17% ...
1	D	227	4% 78% 17% ...
1	E	227	2% 82% 13% ...

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Mol	Chain	Length	Quality of chain
1	F	227	
1	G	227	
1	H	227	
1	I	227	
1	J	227	
1	K	227	
1	L	227	
1	M	227	
1	N	227	
1	O	227	
1	P	227	
1	Q	227	
1	R	227	
1	S	227	
1	T	227	
1	U	227	
1	V	227	
1	W	227	
1	X	227	
1	Y	227	
1	Z	227	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 45004 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 (S)-2-haloacid dehalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	B	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	C	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	D	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	E	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	F	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	G	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	H	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	I	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	J	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	K	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	L	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	M	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	N	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	O	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	P	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	R	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	S	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	T	222	Total	C	N	O	S	0	0	0
			1729	1108	292	322	7			
1	U	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	V	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	W	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	X	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	Y	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			
1	Z	221	Total	C	N	O	S	0	0	0
			1720	1102	290	321	7			

There are 26 discrepancies between the modelled and reference sequences:

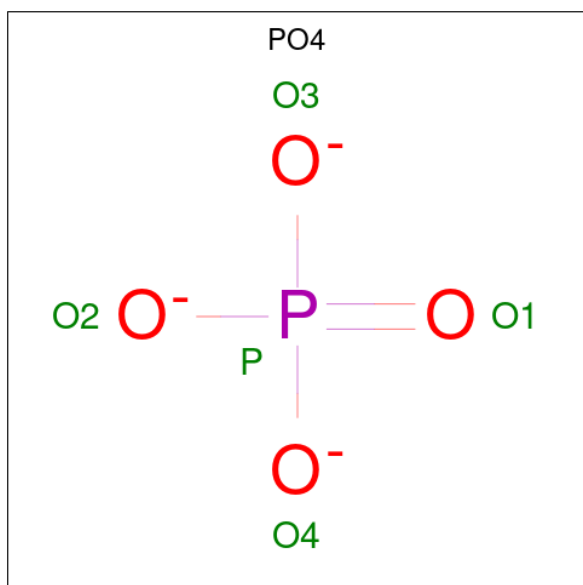
Chain	Residue	Modelled	Actual	Comment	Reference
A	179	ASN	HIS	engineered mutation	UNP G0L7V6
B	179	ASN	HIS	engineered mutation	UNP G0L7V6
C	179	ASN	HIS	engineered mutation	UNP G0L7V6
D	179	ASN	HIS	engineered mutation	UNP G0L7V6
E	179	ASN	HIS	engineered mutation	UNP G0L7V6
F	179	ASN	HIS	engineered mutation	UNP G0L7V6
G	179	ASN	HIS	engineered mutation	UNP G0L7V6
H	179	ASN	HIS	engineered mutation	UNP G0L7V6
I	179	ASN	HIS	engineered mutation	UNP G0L7V6
J	179	ASN	HIS	engineered mutation	UNP G0L7V6
K	179	ASN	HIS	engineered mutation	UNP G0L7V6
L	179	ASN	HIS	engineered mutation	UNP G0L7V6
M	179	ASN	HIS	engineered mutation	UNP G0L7V6
N	179	ASN	HIS	engineered mutation	UNP G0L7V6
O	179	ASN	HIS	engineered mutation	UNP G0L7V6
P	179	ASN	HIS	engineered mutation	UNP G0L7V6
Q	179	ASN	HIS	engineered mutation	UNP G0L7V6
R	179	ASN	HIS	engineered mutation	UNP G0L7V6
S	179	ASN	HIS	engineered mutation	UNP G0L7V6

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Chain	Residue	Modelled	Actual	Comment	Reference
T	179	ASN	HIS	engineered mutation	UNP G0L7V6
U	179	ASN	HIS	engineered mutation	UNP G0L7V6
V	179	ASN	HIS	engineered mutation	UNP G0L7V6
W	179	ASN	HIS	engineered mutation	UNP G0L7V6
X	179	ASN	HIS	engineered mutation	UNP G0L7V6
Y	179	ASN	HIS	engineered mutation	UNP G0L7V6
Z	179	ASN	HIS	engineered mutation	UNP G0L7V6

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	I	1	Total	O	P	0	0
			5	4	1		
2	J	1	Total	O	P	0	0
			5	4	1		
2	K	1	Total	O	P	0	0
			5	4	1		
2	L	1	Total	O	P	0	0
			5	4	1		
2	M	1	Total	O	P	0	0
			5	4	1		
2	N	1	Total	O	P	0	0
			5	4	1		
2	O	1	Total	O	P	0	0
			5	4	1		
2	P	1	Total	O	P	0	0
			5	4	1		
2	Q	1	Total	O	P	0	0
			5	4	1		
2	R	1	Total	O	P	0	0
			5	4	1		
2	S	1	Total	O	P	0	0
			5	4	1		
2	T	1	Total	O	P	0	0
			5	4	1		
2	U	1	Total	O	P	0	0
			5	4	1		
2	V	1	Total	O	P	0	0
			5	4	1		
2	W	1	Total	O	P	0	0
			5	4	1		
2	X	1	Total	O	P	0	0
			5	4	1		
2	Y	1	Total	O	P	0	0
			5	4	1		
2	Z	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	O	0	0
			5	5		

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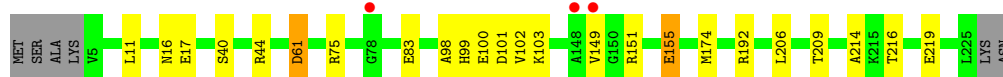
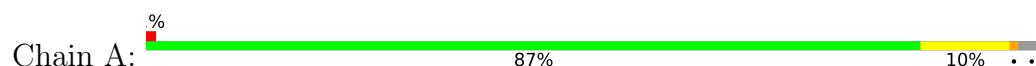
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	O 1	0	0
3	D	2	Total 2	O 2	0	0
3	F	1	Total 1	O 1	0	0
3	G	3	Total 3	O 3	0	0
3	H	4	Total 4	O 4	0	0
3	I	3	Total 3	O 3	0	0
3	J	2	Total 2	O 2	0	0
3	K	6	Total 6	O 6	0	0
3	L	1	Total 1	O 1	0	0
3	M	1	Total 1	O 1	0	0
3	N	2	Total 2	O 2	0	0
3	O	4	Total 4	O 4	0	0
3	P	7	Total 7	O 7	0	0
3	Q	2	Total 2	O 2	0	0
3	R	2	Total 2	O 2	0	0
3	T	2	Total 2	O 2	0	0
3	U	2	Total 2	O 2	0	0
3	W	1	Total 1	O 1	0	0
3	X	2	Total 2	O 2	0	0
3	Y	1	Total 1	O 1	0	0
3	Z	1	Total 1	O 1	0	0

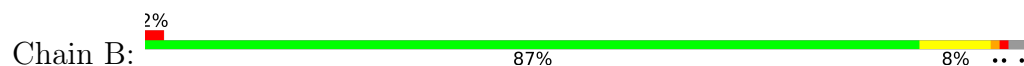
3 Residue-property plots [i](#)

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.

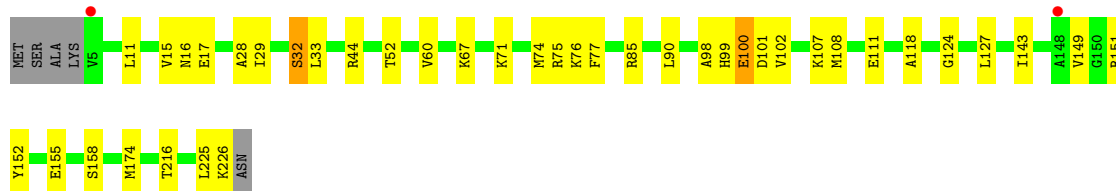
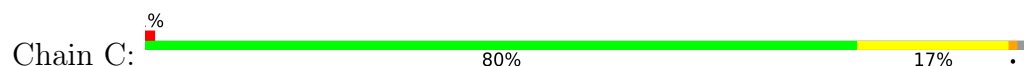
- Molecule 1: (S)-2-haloacid dehalogenase



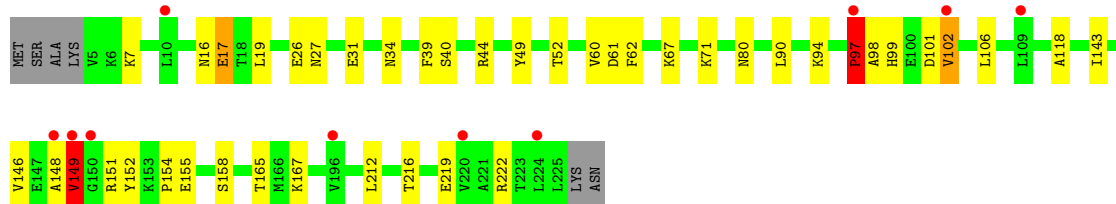
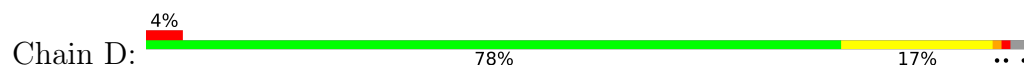
- Molecule 1: (S)-2-haloacid dehalogenase



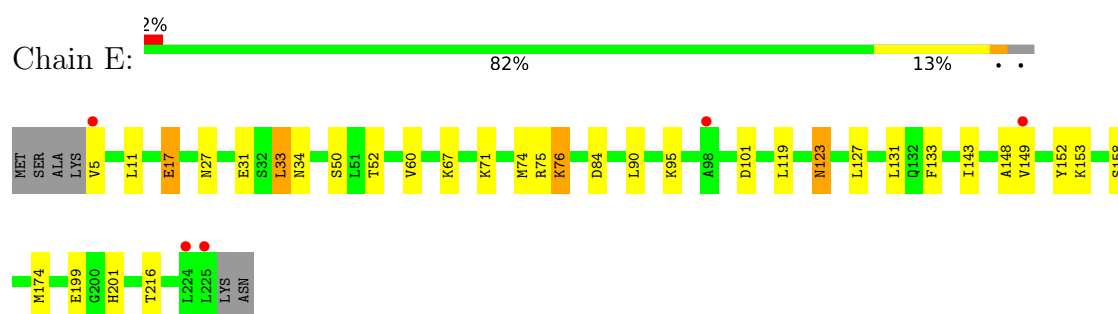
- Molecule 1: (S)-2-haloacid dehalogenase



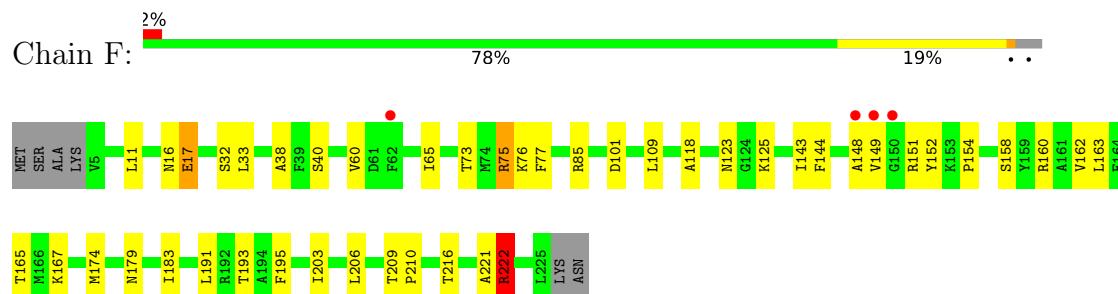
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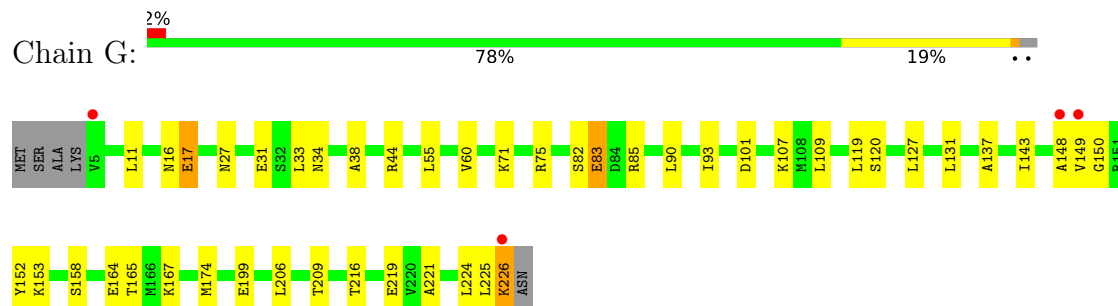
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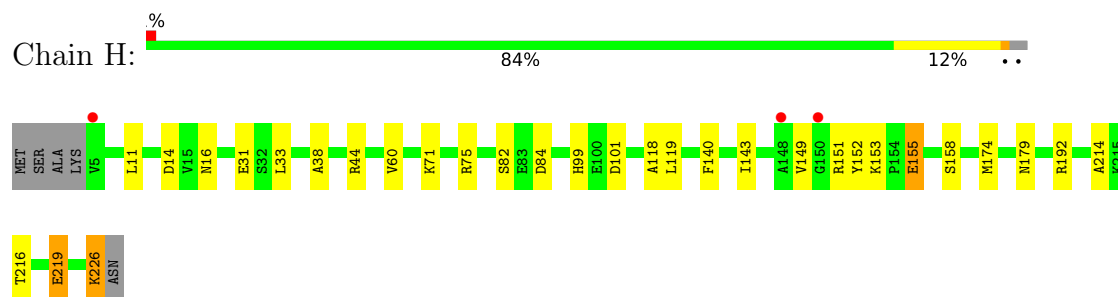
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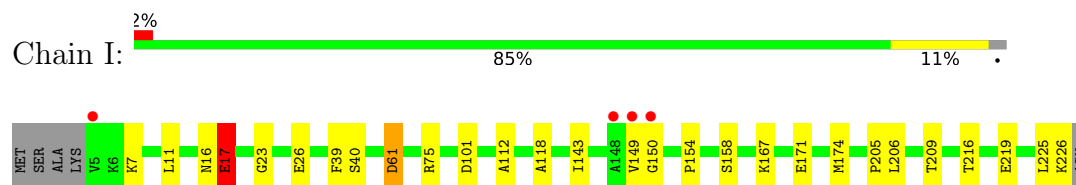
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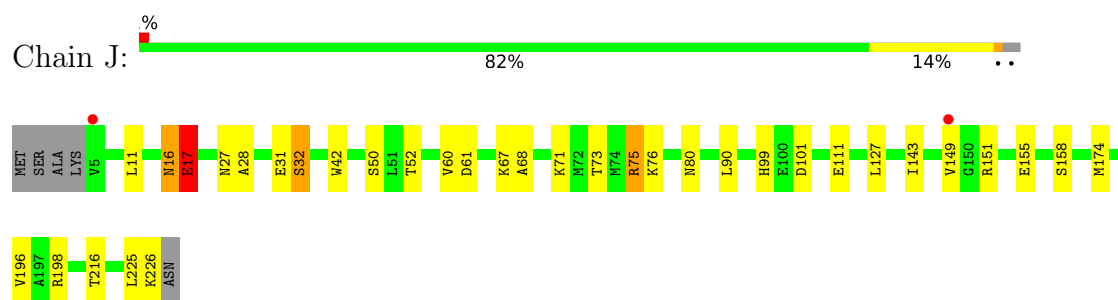
- Molecule 1: (S)-2-haloacid dehalogenase



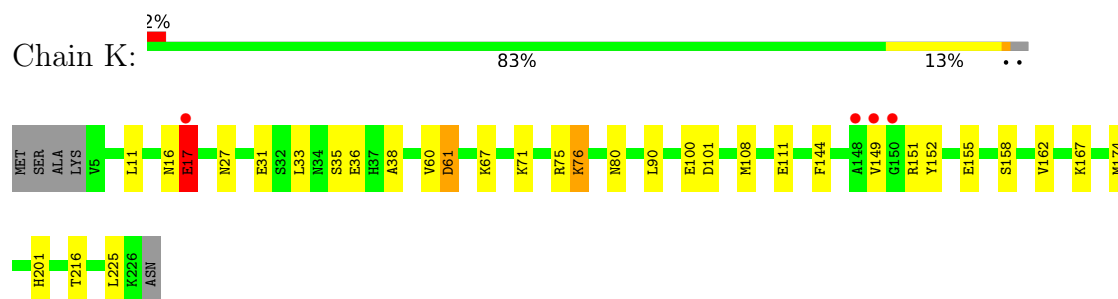
- Molecule 1: (S)-2-haloacid dehalogenase



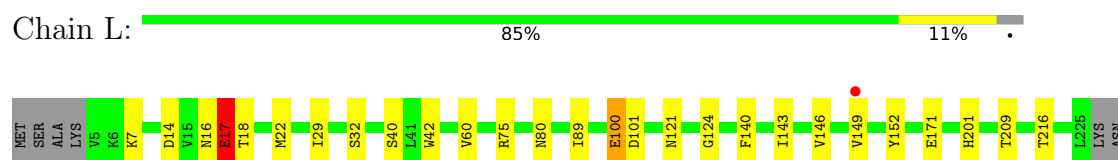
- Molecule 1: (S)-2-haloacid dehalogenase



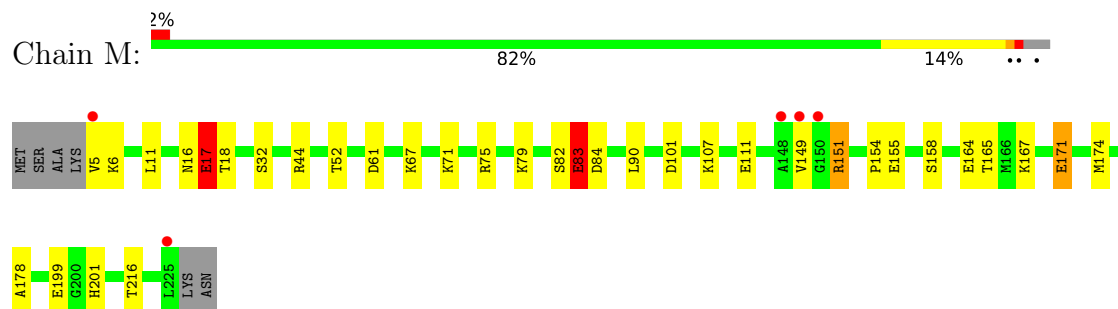
- Molecule 1: (S)-2-haloacid dehalogenase



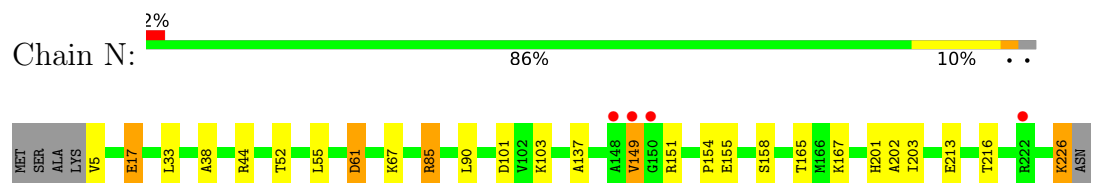
- Molecule 1: (S)-2-haloacid dehalogenase



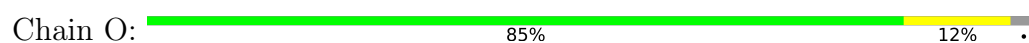
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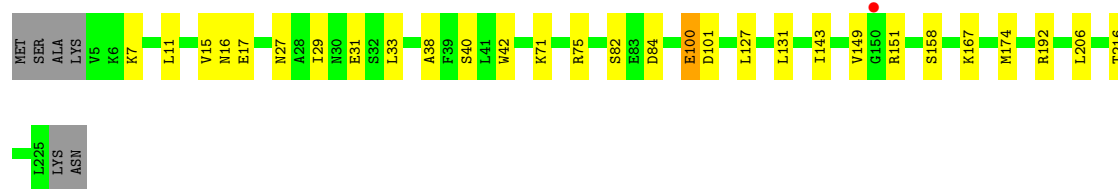


- Molecule 1: (S)-2-haloacid dehalogenase

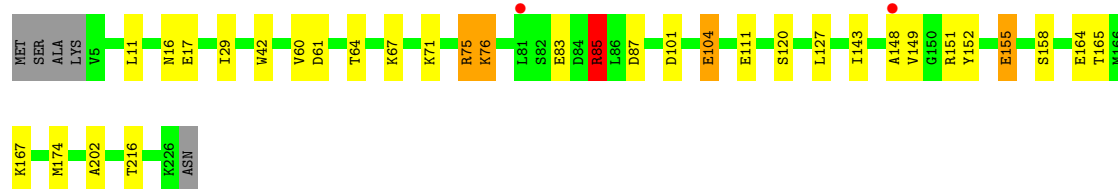
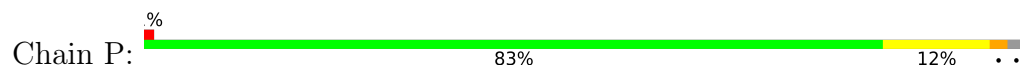


- Molecule 1: (S)-2-haloacid dehalogenase

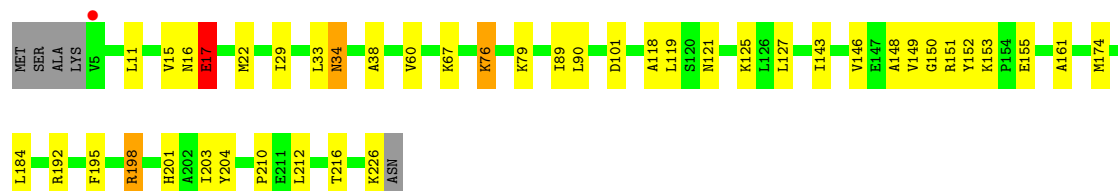
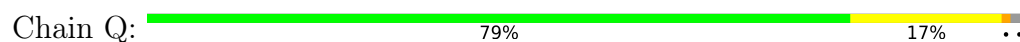




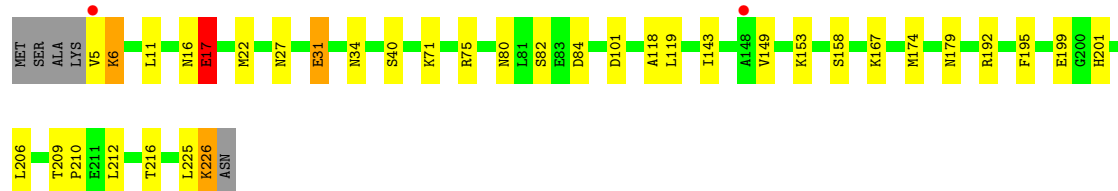
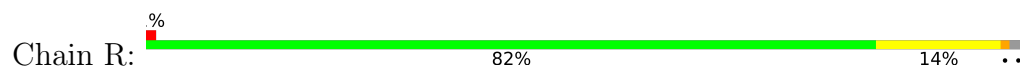
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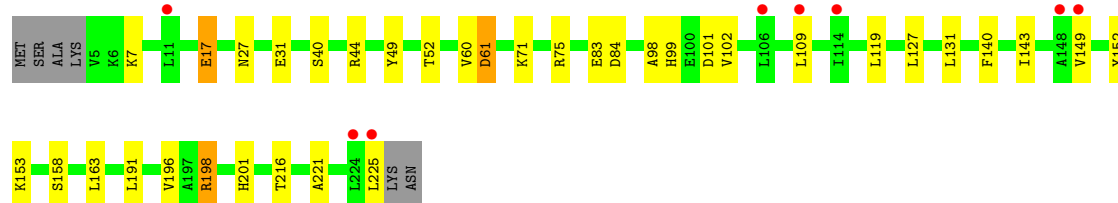
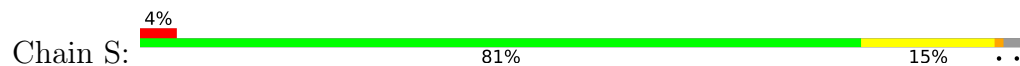
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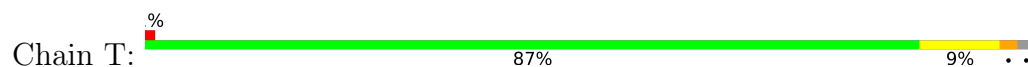
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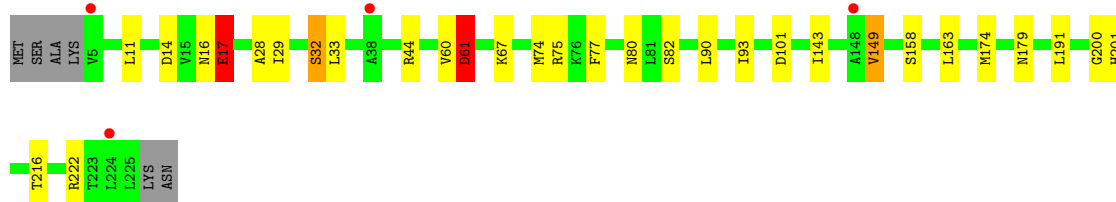
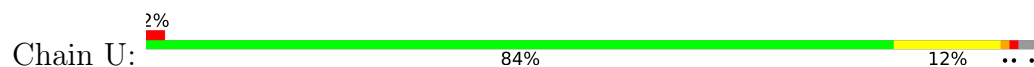
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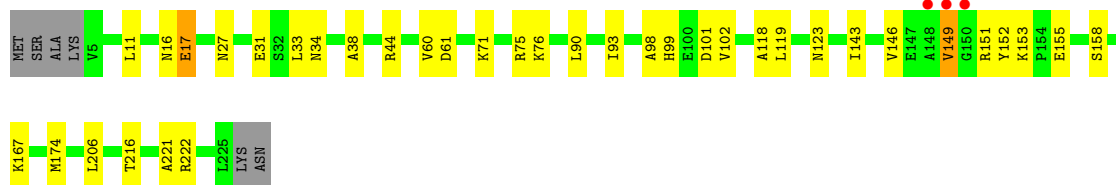
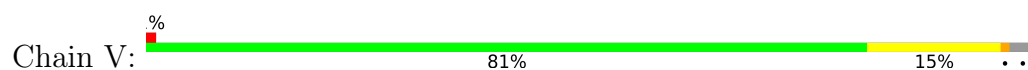
- Molecule 1: (S)-2-haloacid dehalogenase



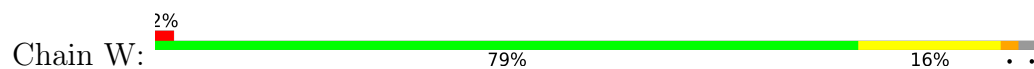
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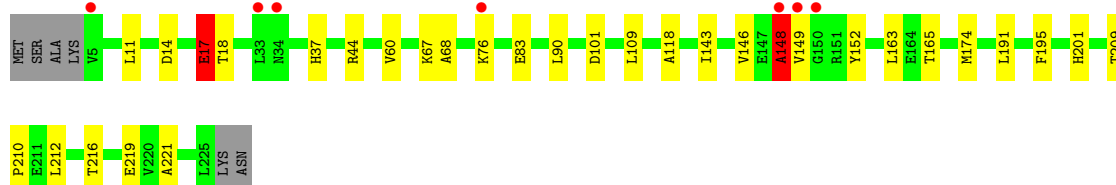
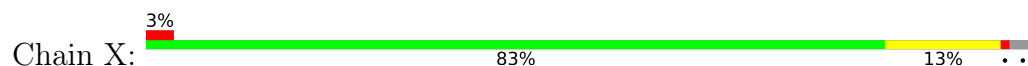
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- Molecule 1: (S)-2-haloacid dehalogenase



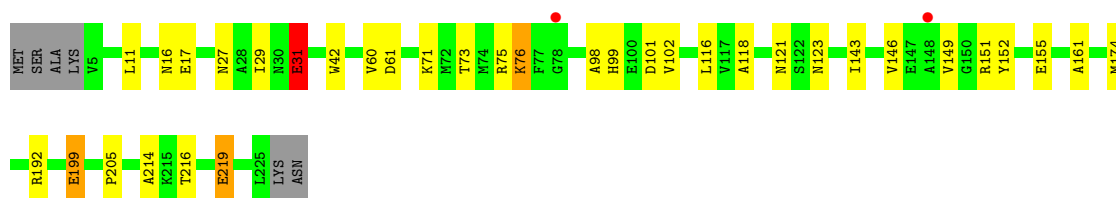
• Molecule 1: (S)-2-haloacid dehalogenase

Chain Y:  % 81% 15% .



• Molecule 1: (S)-2-haloacid dehalogenase

Chain Z:  % 82% 14% . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.17Å 132.79Å 275.70Å 90.00° 92.31° 90.00°	Depositor
Resolution (Å)	49.51 – 2.73 49.51 – 2.73	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.51-2.73) 98.8 (49.51-2.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874, REFMAC 5.8.0267	Depositor
R, R_{free}	0.206 , 0.254 0.210 , 0.254	Depositor DCC
R_{free} test set	7202 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	67.8	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	45004	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.31 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.9379e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PO4

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	0.99	0/1749	1.45	1/2363 (0.0%)
1	B	1.06	2/1749 (0.1%)	1.50	3/2363 (0.1%)
1	C	1.05	1/1758 (0.1%)	1.50	2/2374 (0.1%)
1	D	0.98	0/1749	1.53	4/2363 (0.2%)
1	E	1.03	1/1749 (0.1%)	1.52	2/2363 (0.1%)
1	F	1.10	2/1749 (0.1%)	1.54	8/2363 (0.3%)
1	G	0.98	1/1758 (0.1%)	1.49	4/2374 (0.2%)
1	H	0.99	1/1758 (0.1%)	1.46	4/2374 (0.2%)
1	I	0.98	0/1758	1.44	1/2374 (0.0%)
1	J	0.98	0/1758	1.50	1/2374 (0.0%)
1	K	0.99	1/1758 (0.1%)	1.48	8/2374 (0.3%)
1	L	0.99	1/1749 (0.1%)	1.47	2/2363 (0.1%)
1	M	1.11	5/1749 (0.3%)	1.49	4/2363 (0.2%)
1	N	1.06	1/1758 (0.1%)	1.55	3/2374 (0.1%)
1	O	1.03	1/1749 (0.1%)	1.43	3/2363 (0.1%)
1	P	1.00	1/1758 (0.1%)	1.49	3/2374 (0.1%)
1	Q	1.02	1/1758 (0.1%)	1.54	6/2374 (0.3%)
1	R	1.02	2/1758 (0.1%)	1.49	4/2374 (0.2%)
1	S	0.98	0/1749	1.49	0/2363
1	T	0.98	0/1758	1.46	1/2374 (0.0%)
1	U	0.98	0/1749	1.47	1/2363 (0.0%)
1	V	0.98	0/1749	1.51	3/2363 (0.1%)
1	W	1.11	5/1749 (0.3%)	1.52	5/2363 (0.2%)
1	X	1.05	2/1749 (0.1%)	1.47	3/2363 (0.1%)
1	Y	1.18	5/1749 (0.3%)	1.51	4/2363 (0.2%)
1	Z	1.08	1/1749 (0.1%)	1.50	2/2363 (0.1%)
All	All	1.03	34/45573 (0.1%)	1.49	82/61559 (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	1
1	B	0	1
1	C	0	1
1	F	0	1
1	K	0	1
1	L	0	1
1	M	0	2
1	N	0	1
1	O	0	1
1	P	0	3
1	Q	0	1
1	R	0	1
1	W	0	1
1	X	0	1
1	Z	0	1
All	All	0	18

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	Y	17	GLU	CD-OE2	20.85	1.65	1.25
1	Z	199	GLU	CD-OE2	18.40	1.60	1.25
1	W	5	VAL	C-O	-17.59	0.88	1.23
1	C	100	GLU	CD-OE2	16.43	1.56	1.25
1	N	85	ARG	CZ-NH2	15.92	1.54	1.33

The worst 5 of 82 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	85	ARG	NE-CZ-NH2	20.33	137.50	119.20
1	Q	17	GLU	OE1-CD-OE2	-17.08	81.91	122.90
1	N	85	ARG	NH1-CZ-NH2	-12.07	103.61	119.30
1	M	83	GLU	OE1-CD-OE2	-11.97	94.17	122.90
1	Q	17	GLU	CG-CD-OE1	9.75	140.83	118.40

There are no chirality outliers.

5 of 18 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	GLU	Sidechain
1	B	100	GLU	Sidechain

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Mol	Chain	Res	Type	Group
1	C	100	GLU	Sidechain
1	F	75	ARG	Sidechain
1	K	17	GLU	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	1720	0	1763	19	4
1	B	1720	0	1763	18	4
1	C	1729	0	1776	25	4
1	D	1720	0	1763	25	5
1	E	1720	0	1763	22	0
1	F	1720	0	1763	25	4
1	G	1729	0	1776	29	2
1	H	1729	0	1776	21	0
1	I	1729	0	1776	16	2
1	J	1729	0	1775	32	2
1	K	1729	0	1776	20	1
1	L	1720	0	1763	20	1
1	M	1720	0	1763	30	0
1	N	1729	0	1776	21	1
1	O	1720	0	1763	26	0
1	P	1729	0	1776	36	2
1	Q	1729	0	1776	37	1
1	R	1729	0	1776	33	0
1	S	1720	0	1763	35	1
1	T	1729	0	1776	14	1
1	U	1720	0	1763	25	0
1	V	1720	0	1763	23	3
1	W	1720	0	1763	38	0
1	X	1720	0	1763	18	3
1	Y	1720	0	1763	30	0
1	Z	1720	0	1763	25	1
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	5	0	0	0	0
2	F	5	0	0	0	0
2	G	5	0	0	0	0
2	H	5	0	0	0	0
2	I	5	0	0	0	0
2	J	5	0	0	0	0
2	K	5	0	0	0	0
2	L	5	0	0	0	0
2	M	5	0	0	0	0
2	N	5	0	0	1	0
2	O	5	0	0	0	0
2	P	5	0	0	1	0
2	Q	5	0	0	0	0
2	R	5	0	0	0	0
2	S	5	0	0	0	0
2	T	5	0	0	0	0
2	U	5	0	0	0	0
2	V	5	0	0	0	0
2	W	5	0	0	1	0
2	X	5	0	0	0	0
2	Y	5	0	0	0	0
2	Z	5	0	0	0	0
3	A	5	0	0	3	0
3	B	1	0	0	0	0
3	D	2	0	0	0	0
3	F	1	0	0	0	0
3	G	3	0	0	1	0
3	H	4	0	0	0	0
3	I	3	0	0	1	0
3	J	2	0	0	0	0
3	K	6	0	0	0	0
3	L	1	0	0	0	0
3	M	1	0	0	0	0
3	N	2	0	0	0	0
3	O	4	0	0	0	0
3	P	7	0	0	0	0
3	Q	2	0	0	0	0
3	R	2	0	0	0	0
3	T	2	0	0	0	0
3	U	2	0	0	0	0
3	W	1	0	0	0	0
3	X	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Y	1	0	0	0	0
3	Z	1	0	0	0	0
All	All	45004	0	45980	536	21

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

The worst 5 of 536 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:Y:17:GLU:CD	1:Y:17:GLU:OE2	1.65	1.36
1:J:75:ARG:NH2	1:M:84:ASP:OD1	1.57	1.34
1:H:84:ASP:OD2	1:K:75:ARG:NH2	1.68	1.27
1:W:5:VAL:HG13	1:W:6:LYS:CG	1.70	1.21
1:W:5:VAL:HG12	1:W:6:LYS:HB2	1.26	1.13

The worst 5 of 21 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:111:GLU:OE1	1:S:7:LYS:NZ[2_646]	1.24	0.96
1:A:83:GLU:OE2	1:X:83:GLU:OE2[1_556]	1.47	0.73
1:G:219:GLU:OE1	1:N:103:LYS:NZ[2_646]	1.79	0.41
1:D:97:PRO:CG	1:F:222:ARG:NE[1_455]	1.81	0.39
1:K:111:GLU:OE1	1:L:7:LYS:NZ[1_655]	1.87	0.33

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	219/227 (96%)	215 (98%)	4 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	219/227 (96%)	209 (95%)	8 (4%)	2 (1%)	14	26
1	C	220/227 (97%)	212 (96%)	8 (4%)	0	100	100
1	D	219/227 (96%)	212 (97%)	5 (2%)	2 (1%)	14	26
1	E	219/227 (96%)	214 (98%)	5 (2%)	0	100	100
1	F	219/227 (96%)	212 (97%)	6 (3%)	1 (0%)	24	40
1	G	220/227 (97%)	210 (96%)	9 (4%)	1 (0%)	24	40
1	H	220/227 (97%)	215 (98%)	5 (2%)	0	100	100
1	I	220/227 (97%)	214 (97%)	4 (2%)	2 (1%)	14	26
1	J	220/227 (97%)	212 (96%)	6 (3%)	2 (1%)	14	26
1	K	220/227 (97%)	216 (98%)	4 (2%)	0	100	100
1	L	219/227 (96%)	214 (98%)	5 (2%)	0	100	100
1	M	219/227 (96%)	211 (96%)	7 (3%)	1 (0%)	24	40
1	N	220/227 (97%)	216 (98%)	4 (2%)	0	100	100
1	O	219/227 (96%)	214 (98%)	5 (2%)	0	100	100
1	P	220/227 (97%)	216 (98%)	4 (2%)	0	100	100
1	Q	220/227 (97%)	213 (97%)	6 (3%)	1 (0%)	24	40
1	R	220/227 (97%)	212 (96%)	7 (3%)	1 (0%)	24	40
1	S	219/227 (96%)	213 (97%)	6 (3%)	0	100	100
1	T	220/227 (97%)	212 (96%)	8 (4%)	0	100	100
1	U	219/227 (96%)	209 (95%)	8 (4%)	2 (1%)	14	26
1	V	219/227 (96%)	210 (96%)	7 (3%)	2 (1%)	14	26
1	W	219/227 (96%)	212 (97%)	6 (3%)	1 (0%)	24	40
1	X	219/227 (96%)	217 (99%)	2 (1%)	0	100	100
1	Y	219/227 (96%)	212 (97%)	6 (3%)	1 (0%)	24	40
1	Z	219/227 (96%)	213 (97%)	6 (3%)	0	100	100
All	All	5705/5902 (97%)	5535 (97%)	151 (3%)	19 (0%)	36	54

5 of 19 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	Q	17	GLU
1	U	17	GLU
1	W	6	LYS

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Mol	Chain	Res	Type
1	B	17	GLU
1	D	17	GLU

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	182/187 (97%)	178 (98%)	4 (2%)	45	65
1	B	182/187 (97%)	179 (98%)	3 (2%)	55	72
1	C	183/187 (98%)	181 (99%)	2 (1%)	65	79
1	D	182/187 (97%)	174 (96%)	8 (4%)	25	45
1	E	182/187 (97%)	177 (97%)	5 (3%)	39	61
1	F	182/187 (97%)	177 (97%)	5 (3%)	39	61
1	G	183/187 (98%)	179 (98%)	4 (2%)	45	65
1	H	183/187 (98%)	181 (99%)	2 (1%)	65	79
1	I	183/187 (98%)	179 (98%)	4 (2%)	45	65
1	J	183/187 (98%)	179 (98%)	4 (2%)	45	65
1	K	183/187 (98%)	181 (99%)	2 (1%)	65	79
1	L	182/187 (97%)	180 (99%)	2 (1%)	65	79
1	M	182/187 (97%)	181 (100%)	1 (0%)	81	88
1	N	183/187 (98%)	179 (98%)	4 (2%)	45	65
1	O	182/187 (97%)	181 (100%)	1 (0%)	81	88
1	P	183/187 (98%)	181 (99%)	2 (1%)	65	79
1	Q	183/187 (98%)	181 (99%)	2 (1%)	65	79
1	R	183/187 (98%)	178 (97%)	5 (3%)	39	61
1	S	182/187 (97%)	179 (98%)	3 (2%)	55	72
1	T	183/187 (98%)	176 (96%)	7 (4%)	29	51
1	U	182/187 (97%)	177 (97%)	5 (3%)	39	61
1	V	182/187 (97%)	182 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	W	182/187 (97%)	178 (98%)	4 (2%)	45	65
1	X	182/187 (97%)	179 (98%)	3 (2%)	55	72
1	Y	182/187 (97%)	178 (98%)	4 (2%)	45	65
1	Z	182/187 (97%)	179 (98%)	3 (2%)	55	72
All	All	4743/4862 (98%)	4654 (98%)	89 (2%)	50	68

5 of 89 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	179	ASN
1	U	29	ILE
1	R	212	LEU
1	T	32	SER
1	W	7	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 31 such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	HIS
1	U	80	ASN
1	L	179	ASN
1	Z	16	ASN
1	R	172	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

26 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	PO4	S	301	-	4,4,4	0.76	0	6,6,6	0.53	0
2	PO4	Q	301	-	4,4,4	0.82	0	6,6,6	0.45	0
2	PO4	T	301	-	4,4,4	0.87	0	6,6,6	0.41	0
2	PO4	X	301	-	4,4,4	0.96	0	6,6,6	0.43	0
2	PO4	A	301	-	4,4,4	0.63	0	6,6,6	0.55	0
2	PO4	E	301	-	4,4,4	0.65	0	6,6,6	0.48	0
2	PO4	Z	301	-	4,4,4	0.76	0	6,6,6	0.47	0
2	PO4	O	301	-	4,4,4	0.75	0	6,6,6	0.56	0
2	PO4	L	301	-	4,4,4	0.77	0	6,6,6	0.49	0
2	PO4	C	301	-	4,4,4	0.66	0	6,6,6	0.48	0
2	PO4	N	301	-	4,4,4	1.02	0	6,6,6	0.40	0
2	PO4	V	301	-	4,4,4	0.62	0	6,6,6	0.54	0
2	PO4	B	301	-	4,4,4	0.75	0	6,6,6	0.51	0
2	PO4	D	301	-	4,4,4	0.58	0	6,6,6	0.58	0
2	PO4	M	301	-	4,4,4	0.83	0	6,6,6	0.44	0
2	PO4	U	301	-	4,4,4	0.72	0	6,6,6	0.45	0
2	PO4	K	301	-	4,4,4	0.74	0	6,6,6	0.51	0
2	PO4	G	301	-	4,4,4	1.03	0	6,6,6	0.37	0
2	PO4	R	301	-	4,4,4	0.69	0	6,6,6	0.53	0
2	PO4	F	301	-	4,4,4	0.68	0	6,6,6	0.49	0
2	PO4	P	301	-	4,4,4	0.74	0	6,6,6	0.58	0
2	PO4	Y	301	-	4,4,4	0.62	0	6,6,6	0.52	0
2	PO4	J	301	-	4,4,4	0.98	0	6,6,6	0.42	0
2	PO4	W	301	-	4,4,4	1.05	0	6,6,6	0.48	0
2	PO4	H	301	-	4,4,4	0.73	0	6,6,6	0.54	0
2	PO4	I	301	-	4,4,4	0.55	0	6,6,6	0.53	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	N	301	PO4	1	0
2	P	301	PO4	1	0
2	W	301	PO4	1	0

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	221/227 (97%)	0.09	3 (1%) 73 72	61, 84, 123, 138	0
1	B	221/227 (97%)	0.06	4 (1%) 67 64	64, 85, 116, 135	0
1	C	222/227 (97%)	0.09	2 (0%) 81 81	67, 92, 120, 139	0
1	D	221/227 (97%)	0.53	10 (4%) 38 37	72, 108, 144, 190	0
1	E	221/227 (97%)	0.51	5 (2%) 61 58	72, 100, 136, 168	0
1	F	221/227 (97%)	0.38	4 (1%) 67 64	72, 89, 121, 172	0
1	G	222/227 (97%)	-0.04	4 (1%) 67 64	59, 78, 110, 147	0
1	H	222/227 (97%)	-0.06	3 (1%) 73 72	58, 76, 106, 157	0
1	I	222/227 (97%)	0.00	4 (1%) 67 64	56, 72, 99, 123	0
1	J	222/227 (97%)	-0.06	2 (0%) 81 81	59, 76, 112, 139	0
1	K	222/227 (97%)	-0.12	4 (1%) 67 64	54, 72, 102, 121	0
1	L	221/227 (97%)	-0.06	1 (0%) 87 87	54, 73, 98, 117	0
1	M	221/227 (97%)	0.09	5 (2%) 61 58	61, 83, 118, 164	0
1	N	222/227 (97%)	0.01	4 (1%) 67 64	60, 75, 108, 132	0
1	O	221/227 (97%)	-0.21	1 (0%) 87 87	52, 66, 99, 134	0
1	P	222/227 (97%)	-0.16	2 (0%) 81 81	52, 68, 98, 117	0
1	Q	222/227 (97%)	0.08	1 (0%) 87 87	56, 83, 115, 139	0
1	R	222/227 (97%)	0.08	2 (0%) 81 81	63, 90, 121, 147	0
1	S	221/227 (97%)	0.56	8 (3%) 46 44	66, 97, 126, 153	0
1	T	222/227 (97%)	0.17	3 (1%) 73 72	69, 93, 125, 144	0
1	U	221/227 (97%)	0.20	4 (1%) 67 64	65, 90, 118, 145	0
1	V	221/227 (97%)	0.23	3 (1%) 73 72	65, 96, 129, 140	0
1	W	221/227 (97%)	0.34	5 (2%) 61 58	68, 97, 131, 161	0
1	X	221/227 (97%)	0.24	7 (3%) 50 48	68, 92, 124, 154	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Y	221/227 (97%)	0.26	2 (0%) 81 81	73, 100, 131, 142	0
1	Z	221/227 (97%)	0.24	2 (0%) 81 81	67, 98, 135, 157	0
All	All	5757/5902 (97%)	0.13	95 (1%) 69 66	52, 85, 124, 190	0

The worst 5 of 95 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	ALA	4.9
1	P	148	ALA	4.4
1	X	148	ALA	4.2
1	B	148	ALA	4.0
1	V	148	ALA	3.9

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PO4	F	301	5/5	0.44	0.15	126,153,162,167	0
2	PO4	E	301	5/5	0.76	0.12	130,143,144,153	0
2	PO4	W	301	5/5	0.79	0.14	91,97,109,119	0
2	PO4	Q	301	5/5	0.81	0.11	101,139,142,144	0
2	PO4	D	301	5/5	0.83	0.10	92,104,111,127	0
2	PO4	Z	301	5/5	0.84	0.10	110,126,133,139	0
2	PO4	Y	301	5/5	0.85	0.11	126,137,146,154	0
2	PO4	I	301	5/5	0.86	0.14	73,98,118,124	0
2	PO4	M	301	5/5	0.87	0.10	97,123,130,140	0
2	PO4	B	301	5/5	0.89	0.10	94,99,103,107	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	N	301	5/5	0.89	0.09	98,105,112,122	0
2	PO4	C	301	5/5	0.90	0.10	101,120,122,128	0
2	PO4	J	301	5/5	0.90	0.11	107,114,117,123	0
2	PO4	X	301	5/5	0.91	0.13	90,99,107,108	0
2	PO4	K	301	5/5	0.91	0.14	77,106,113,119	0
2	PO4	P	301	5/5	0.91	0.14	76,77,92,97	0
2	PO4	L	301	5/5	0.92	0.14	93,103,111,112	0
2	PO4	T	301	5/5	0.92	0.10	101,103,106,108	0
2	PO4	G	301	5/5	0.93	0.15	77,82,88,102	0
2	PO4	V	301	5/5	0.93	0.09	82,82,92,98	0
2	PO4	O	301	5/5	0.94	0.12	76,82,94,96	0
2	PO4	R	301	5/5	0.94	0.09	99,100,109,114	0
2	PO4	S	301	5/5	0.94	0.08	96,99,105,109	0
2	PO4	A	301	5/5	0.94	0.09	76,89,92,99	0
2	PO4	U	301	5/5	0.94	0.12	88,94,102,103	0
2	PO4	H	301	5/5	0.96	0.07	74,86,86,89	0

6.5 Other polymers [i](#)

There are no such residues in this entry.