



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:44 PM UTC

PDB ID : 9H0Z / pdb_00009h0z
Title : Crystal structure of TTL[Nle], thermophilic lipase TTL from Thermoanaerobacter thermohydrosulfuricus containing non-canonical amino acid Nle at the position of Met
Authors : Hromic-Jahjefendic, A.; Pavkov-Keller, T.; Wiltshi, B.; Gruber, K.
Deposited on : 2024-10-09
Resolution : 2.79 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
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

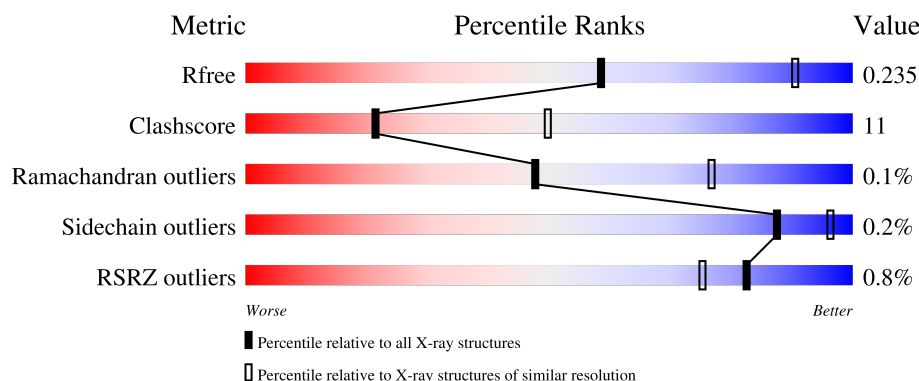
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	267	 74% 22% .
1	B	267	 68% 28% .
1	C	267	 71% 22% 7%
1	D	267	 2% 68% 25% 7%
1	E	267	 74% 19% 7%

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Mol	Chain	Length	Quality of chain
1	F	267	<div><div>%</div><div><div></div><div>71%</div><div>24%</div><div>5%</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12131 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 Serine aminopeptidase S33 domain-containing protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	0	1	0
			2046	1322	334	390			
1	B	257	Total	C	N	O	0	0	0
			2037	1317	333	387			
1	C	248	Total	C	N	O	0	0	0
			1960	1267	322	371			
1	D	247	Total	C	N	O	0	0	0
			1957	1266	320	371			
1	E	249	Total	C	N	O	0	1	0
			1979	1279	324	376			
1	F	253	Total	C	N	O	0	0	0
			2000	1292	327	381			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
A	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
A	260	GLY	-	expression tag	UNP A0A1G7VV58
A	261	SER	-	expression tag	UNP A0A1G7VV58
A	262	HIS	-	expression tag	UNP A0A1G7VV58
A	263	HIS	-	expression tag	UNP A0A1G7VV58
A	264	HIS	-	expression tag	UNP A0A1G7VV58

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Chain	Residue	Modelled	Actual	Comment	Reference
A	265	HIS	-	expression tag	UNP A0A1G7VV58
A	266	HIS	-	expression tag	UNP A0A1G7VV58
A	267	HIS	-	expression tag	UNP A0A1G7VV58
B	1	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
B	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
B	260	GLY	-	expression tag	UNP A0A1G7VV58
B	261	SER	-	expression tag	UNP A0A1G7VV58
B	262	HIS	-	expression tag	UNP A0A1G7VV58
B	263	HIS	-	expression tag	UNP A0A1G7VV58
B	264	HIS	-	expression tag	UNP A0A1G7VV58
B	265	HIS	-	expression tag	UNP A0A1G7VV58
B	266	HIS	-	expression tag	UNP A0A1G7VV58
B	267	HIS	-	expression tag	UNP A0A1G7VV58
C	1	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
C	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
C	260	GLY	-	expression tag	UNP A0A1G7VV58
C	261	SER	-	expression tag	UNP A0A1G7VV58
C	262	HIS	-	expression tag	UNP A0A1G7VV58
C	263	HIS	-	expression tag	UNP A0A1G7VV58
C	264	HIS	-	expression tag	UNP A0A1G7VV58
C	265	HIS	-	expression tag	UNP A0A1G7VV58
C	266	HIS	-	expression tag	UNP A0A1G7VV58

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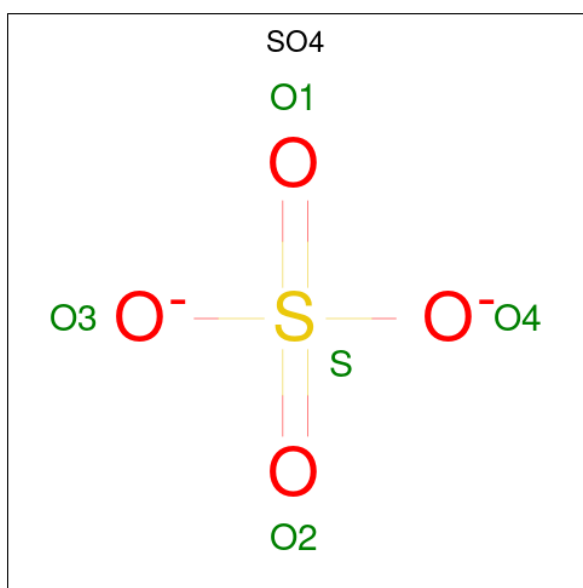
Chain	Residue	Modelled	Actual	Comment	Reference
C	267	HIS	-	expression tag	UNP A0A1G7VV58
D	1	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
D	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
D	260	GLY	-	expression tag	UNP A0A1G7VV58
D	261	SER	-	expression tag	UNP A0A1G7VV58
D	262	HIS	-	expression tag	UNP A0A1G7VV58
D	263	HIS	-	expression tag	UNP A0A1G7VV58
D	264	HIS	-	expression tag	UNP A0A1G7VV58
D	265	HIS	-	expression tag	UNP A0A1G7VV58
D	266	HIS	-	expression tag	UNP A0A1G7VV58
D	267	HIS	-	expression tag	UNP A0A1G7VV58
E	1	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
E	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
E	260	GLY	-	expression tag	UNP A0A1G7VV58
E	261	SER	-	expression tag	UNP A0A1G7VV58
E	262	HIS	-	expression tag	UNP A0A1G7VV58
E	263	HIS	-	expression tag	UNP A0A1G7VV58
E	264	HIS	-	expression tag	UNP A0A1G7VV58
E	265	HIS	-	expression tag	UNP A0A1G7VV58
E	266	HIS	-	expression tag	UNP A0A1G7VV58
E	267	HIS	-	expression tag	UNP A0A1G7VV58
F	1	NLE	MET	engineered mutation	UNP A0A1G7VV58

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Chain	Residue	Modelled	Actual	Comment	Reference
F	17	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	18	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	30	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	33	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	51	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	80	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	114	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	142	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	147	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	148	ASN	HIS	engineered mutation	UNP A0A1G7VV58
F	158	NLE	MET	engineered mutation	UNP A0A1G7VV58
F	260	GLY	-	expression tag	UNP A0A1G7VV58
F	261	SER	-	expression tag	UNP A0A1G7VV58
F	262	HIS	-	expression tag	UNP A0A1G7VV58
F	263	HIS	-	expression tag	UNP A0A1G7VV58
F	264	HIS	-	expression tag	UNP A0A1G7VV58
F	265	HIS	-	expression tag	UNP A0A1G7VV58
F	266	HIS	-	expression tag	UNP A0A1G7VV58
F	267	HIS	-	expression tag	UNP A0A1G7VV58

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

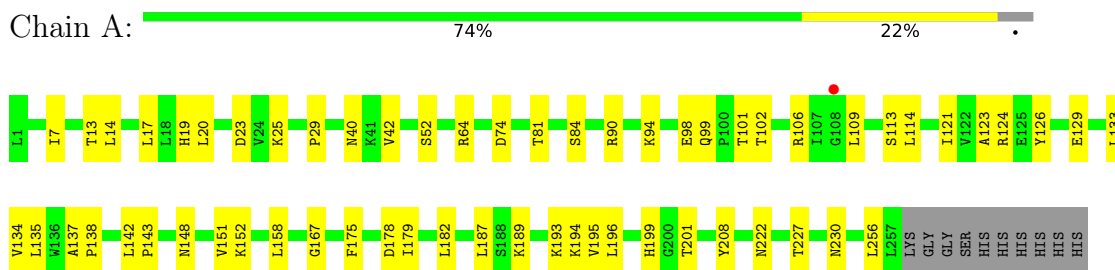
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		
3	B	14	Total	O	0	0
			14	14		
3	C	26	Total	O	0	0
			26	26		
3	D	22	Total	O	0	0
			22	22		
3	E	23	Total	O	0	0
			23	23		
3	F	20	Total	O	0	0
			20	20		

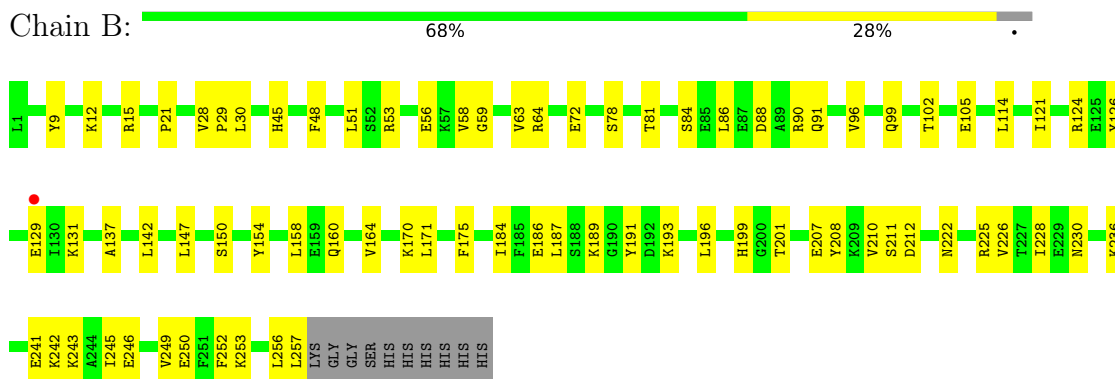
3 Residue-property plots

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.

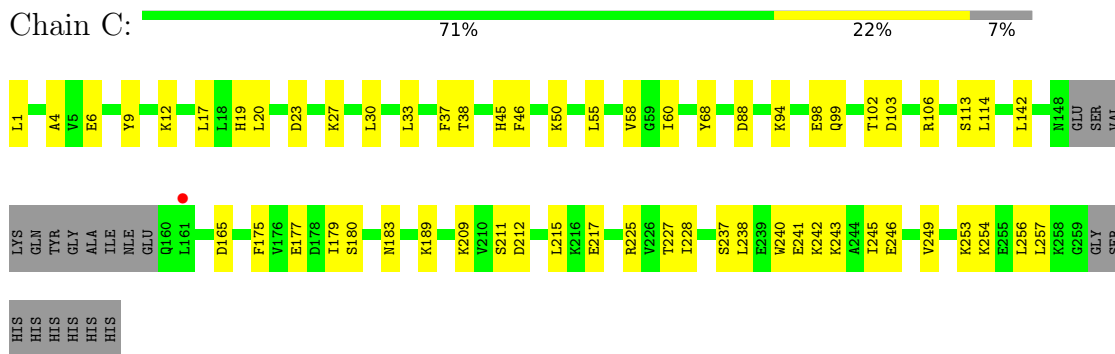
- Molecule 1: Serine aminopeptidase S33 domain-containing protein



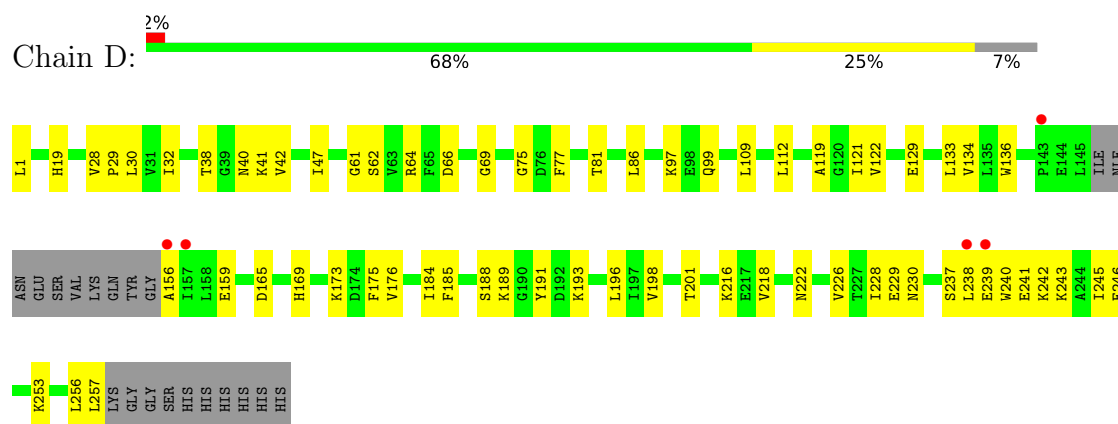
- Molecule 1: Serine aminopeptidase S33 domain-containing protein



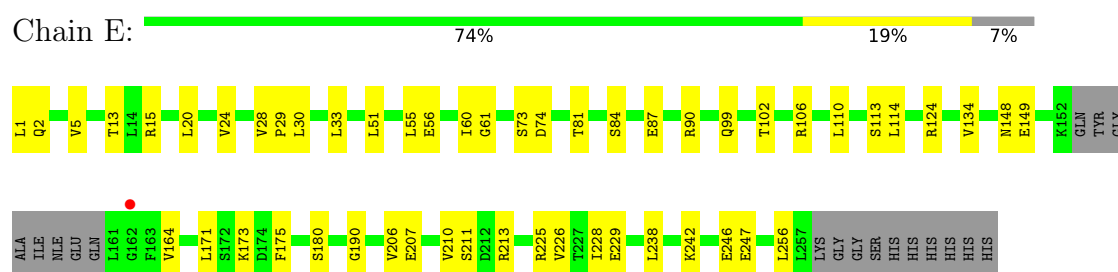
- Molecule 1: Serine aminopeptidase S33 domain-containing protein



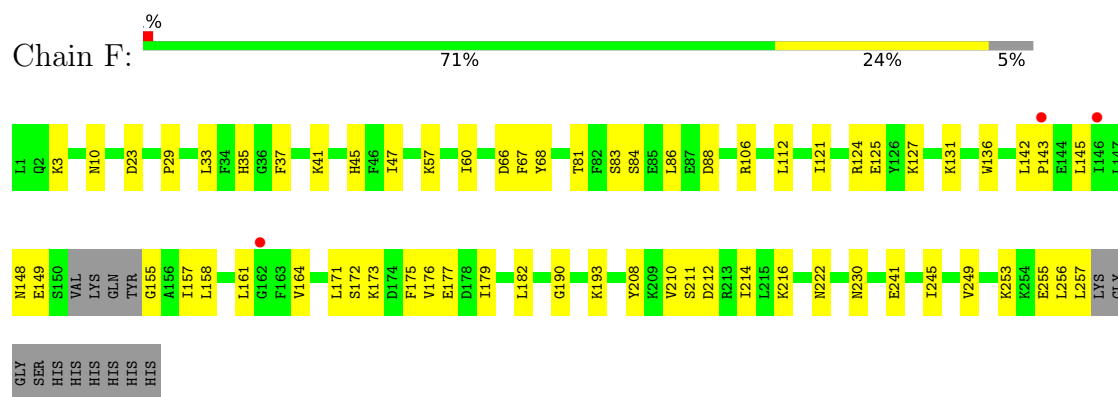
- Molecule 1: Serine aminopeptidase S33 domain-containing protein



- Molecule 1: Serine aminopeptidase S33 domain-containing protein



- Molecule 1: Serine aminopeptidase S33 domain-containing protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.90Å 76.30Å 187.79Å 90.00° 93.52° 90.00°	Depositor
Resolution (Å)	54.89 – 2.79 54.89 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.2 (54.89-2.79) 61.2 (54.89-2.79)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.214 , 0.263 (Not available) , 0.235	Depositor DCC
R_{free} test set	2267 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12131	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.66% 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: NLE, SO4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.14	0/1987	0.34	0/2658
1	B	0.12	0/1978	0.34	0/2646
1	C	0.13	0/1908	0.34	0/2552
1	D	0.21	0/1905	0.40	0/2548
1	E	0.14	0/1927	0.35	0/2577
1	F	0.14	0/1939	0.35	0/2592
All	All	0.15	0/11644	0.35	0/15573

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2046	0	2058	40	1
1	B	2037	0	2053	56	0
1	C	1960	0	1972	42	1
1	D	1957	0	1972	50	0
1	E	1979	0	1996	34	0
1	F	2000	0	2014	48	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	17	0	0	6	0
3	B	14	0	0	2	0
3	C	26	0	0	4	0
3	D	22	0	0	5	0
3	E	23	0	0	7	0
3	F	20	0	0	9	0
All	All	12131	0	12065	262	1

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 11.

All (262) 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:F:177:GLU:OE2	3:F:401:HOH:O	1.86	0.93
1:D:189:LYS:HG2	1:D:218:VAL:HA	1.51	0.92
1:B:150:SER:HA	1:B:154:TYR:HB2	1.54	0.89
1:E:246:GLU:O	3:E:402:HOH:O	1.92	0.87
1:F:23:ASP:OD1	3:F:402:HOH:O	1.98	0.81
1:F:88:ASP:OD2	3:F:403:HOH:O	2.00	0.79
1:C:33:NLE:O	3:C:401:HOH:O	2.00	0.79
1:F:145:LEU:O	1:F:149:GLU:OE1	2.01	0.79
1:E:229:GLU:O	3:E:403:HOH:O	2.03	0.77
1:F:230:ASN:O	3:F:404:HOH:O	2.04	0.76
1:B:196:LEU:HD11	1:B:226:VAL:HG23	1.68	0.73
1:A:113:SER:OG	1:A:114:NLE:N	2.22	0.71
1:B:29:PRO:HG3	1:B:257:LEU:HD21	1.70	0.71
1:D:1:NLE:N	3:D:404:HOH:O	2.23	0.71
1:D:228:ILE:HG12	1:D:243:LYS:HD3	1.74	0.70
1:B:90:ARG:NH2	1:F:23:ASP:O	2.26	0.69
1:F:10:ASN:ND2	3:F:408:HOH:O	2.26	0.69
1:A:109:LEU:HB2	1:A:133:LEU:HD12	1.75	0.68
1:C:183:ASN:ND2	3:C:404:HOH:O	2.27	0.68
1:D:246:GLU:OE2	3:D:401:HOH:O	2.11	0.67
1:F:41:LYS:NZ	3:F:405:HOH:O	2.17	0.66
1:F:148:ASN:HB2	1:F:149:GLU:OE1	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:207:GLU:O	1:E:210:VAL:HG12	1.96	0.64
1:A:194:LYS:NZ	3:A:408:HOH:O	2.30	0.64
1:D:30:NLE:HD3	1:D:61:GLY:HA3	1.78	0.64
1:A:134:VAL:HG22	1:A:196:LEU:HB3	1.80	0.64
1:A:152:LYS:N	3:A:402:HOH:O	2.23	0.64
1:B:45:HIS:ND1	1:B:241:GLU:OE2	2.31	0.63
1:F:158:NLE:HD2	1:F:176:VAL:HG11	1.80	0.63
1:F:155:GLY:N	3:F:410:HOH:O	2.31	0.63
1:D:134:VAL:HG22	1:D:196:LEU:HB3	1.81	0.62
1:F:45:HIS:ND1	1:F:241:GLU:OE2	2.29	0.62
1:A:151:VAL:N	3:A:402:HOH:O	2.30	0.62
1:C:254:LYS:NZ	3:C:405:HOH:O	2.32	0.61
1:E:210:VAL:HA	1:E:213:ARG:HD3	1.82	0.61
1:C:45:HIS:O	1:C:50:LYS:NZ	2.28	0.61
1:B:21:PRO:HG3	1:B:28:VAL:HG21	1.82	0.60
1:D:28:VAL:HG22	1:D:29:PRO:HD2	1.84	0.60
1:A:17:NLE:HE3	1:B:15:ARG:HD3	1.84	0.60
1:B:186:GLU:OE2	1:F:57:LYS:NZ	2.34	0.60
1:C:237:SER:HG	1:C:240:TRP:H	1.49	0.59
1:F:68:TYR:HB3	3:F:403:HOH:O	2.01	0.59
1:A:123:ALA:HB2	1:A:133:LEU:HD13	1.84	0.59
1:C:215:LEU:HD11	1:C:225:ARG:HB2	1.85	0.58
1:D:156:ALA:HA	1:D:159:GLU:HG2	1.85	0.58
1:C:27:LYS:HB3	1:C:103:ASP:HB3	1.84	0.58
1:B:201:THR:HB	1:B:230:ASN:H	1.67	0.58
1:C:99:GLN:HB2	1:C:102:THR:HG23	1.84	0.58
1:A:148:ASN:O	3:A:402:HOH:O	2.17	0.58
1:B:252:PHE:O	1:B:256:LEU:HB2	2.04	0.58
1:D:97:LYS:NZ	1:D:129:GLU:OE2	2.36	0.58
1:F:112:LEU:HD12	1:F:136:TRP:HB3	1.86	0.58
1:A:126:TYR:HB3	1:A:129:GLU:HB2	1.84	0.58
1:A:81:THR:HA	1:A:175:PHE:HD1	1.69	0.57
1:E:106:ARG:NH1	1:E:256:LEU:O	2.30	0.57
1:D:239:GLU:HA	1:D:242:LYS:CB	2.34	0.57
1:F:47:ILE:HD12	1:F:241:GLU:HG3	1.86	0.57
1:B:250:GLU:HA	1:B:253:LYS:HG2	1.84	0.57
1:C:237:SER:HG	1:C:240:TRP:CG	2.23	0.57
1:B:81:THR:H	1:B:84:SER:HB3	1.69	0.57
1:D:47:ILE:HG13	1:D:245:ILE:HD11	1.86	0.57
1:D:198:VAL:HG22	1:D:226:VAL:HG23	1.85	0.57
1:F:253:LYS:HG3	1:F:257:LEU:HD12	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:211:SER:OG	1:E:225:ARG:NH1	2.38	0.56
1:B:58:VAL:HG11	1:B:253:LYS:HB3	1.88	0.56
1:F:179:ILE:HA	1:F:182:LEU:HD12	1.88	0.56
1:E:226:VAL:HG21	1:E:247:GLU:HG2	1.88	0.55
1:C:189:LYS:NZ	1:C:217:GLU:O	2.37	0.55
1:F:33:NLE:O	3:F:406:HOH:O	2.18	0.55
1:C:45:HIS:ND1	1:C:241:GLU:OE2	2.34	0.55
1:C:106:ARG:HD3	1:C:256:LEU:O	2.06	0.55
1:A:199:HIS:CD2	1:A:208:TYR:HA	2.41	0.55
1:B:208:TYR:O	1:B:211:SER:OG	2.22	0.54
1:F:148:ASN:CB	1:F:149:GLU:OE1	2.54	0.54
1:A:178:ASP:OD1	3:A:403:HOH:O	2.18	0.54
1:F:193:LYS:C	1:F:222:ASN:HD21	2.16	0.54
1:A:208:TYR:HH	1:A:227:THR:HG1	1.47	0.54
1:C:245:ILE:O	1:C:249:VAL:HG23	2.07	0.54
1:B:207:GLU:O	1:B:210:VAL:HG12	2.07	0.54
1:C:55:LEU:HD23	1:C:249:VAL:HG13	1.88	0.54
1:F:158:NLE:HE2	1:F:164:VAL:HB	1.89	0.54
1:D:185:PHE:O	1:D:189:LYS:HG3	2.08	0.53
1:B:96:VAL:HG12	1:B:102:THR:HG21	1.90	0.53
1:B:199:HIS:CD2	1:B:208:TYR:HA	2.44	0.53
1:B:245:ILE:O	1:B:249:VAL:HG23	2.09	0.53
1:E:113:SER:OG	1:E:114:NLE:N	2.40	0.53
1:B:78:SER:HB3	1:B:170:LYS:HE2	1.91	0.52
1:D:239:GLU:HA	1:D:242:LYS:HB2	1.91	0.52
1:E:13:THR:HB	1:E:74:ASP:OD2	2.10	0.52
1:A:179:ILE:HA	1:A:182:LEU:HD12	1.92	0.52
1:A:124:ARG:HG3	1:A:189:LYS:O	2.09	0.52
1:B:12:LYS:NZ	3:B:403:HOH:O	2.28	0.52
1:C:113:SER:OG	1:C:114:NLE:N	2.42	0.52
1:D:242:LYS:HG3	1:D:246:GLU:OE1	2.09	0.52
1:D:41:LYS:N	1:D:66:ASP:OD1	2.35	0.52
1:D:41:LYS:HB2	1:D:64:ARG:HB3	1.91	0.52
1:A:81:THR:HG23	1:A:84:SER:H	1.74	0.51
1:E:164:VAL:HG13	1:E:171:LEU:HB3	1.92	0.51
1:C:242:LYS:O	1:C:246:GLU:HB2	2.10	0.51
1:B:164:VAL:HG12	1:B:171:LEU:HB3	1.91	0.51
1:A:106:ARG:HG2	1:A:256:LEU:O	2.10	0.51
1:B:252:PHE:HD1	1:B:256:LEU:HD13	1.76	0.50
1:C:20:LEU:HD21	1:C:30:NLE:HE1	1.92	0.50
1:F:81:THR:HA	1:F:175:PHE:CD2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:PRO:HB3	1:B:59:GLY:HA2	1.94	0.50
1:F:60:ILE:HD13	1:F:256:LEU:HD12	1.92	0.50
1:F:172:SER:O	1:F:176:VAL:HG13	2.10	0.50
1:B:184:ILE:N	3:B:406:HOH:O	2.37	0.50
1:B:212:ASP:OD1	1:B:225:ARG:NH1	2.45	0.50
1:F:164:VAL:HG12	1:F:171:LEU:HB3	1.92	0.50
1:B:86:LEU:HD21	1:B:121:ILE:HG21	1.92	0.50
1:B:114:NLE:HD3	1:B:142:NLE:HE3	1.93	0.50
1:D:257:LEU:O	3:D:402:HOH:O	2.20	0.50
1:C:46:PHE:HE1	1:D:77:PHE:HE1	1.60	0.50
1:C:227:THR:O	1:C:243:LYS:NZ	2.45	0.50
1:D:239:GLU:HA	1:D:242:LYS:HB3	1.94	0.50
1:F:157:ILE:HD12	1:F:157:ILE:H	1.76	0.49
1:C:1:NLE:O	1:C:19:HIS:HA	2.12	0.49
1:C:4:ALA:HB2	1:C:17:NLE:HE2	1.94	0.49
1:C:94:LYS:O	1:C:98:GLU:HG3	2.13	0.49
1:E:247:GLU:OE1	3:E:404:HOH:O	2.20	0.49
1:F:245:ILE:O	1:F:249:VAL:HG23	2.13	0.49
1:B:99:GLN:HB2	1:B:102:THR:OG1	2.12	0.49
1:D:237:SER:HB3	1:D:240:TRP:CD2	2.47	0.49
1:B:51:NLE:HD2	1:B:245:ILE:HG23	1.95	0.48
1:B:81:THR:HA	1:B:175:PHE:HD2	1.78	0.48
1:C:58:VAL:HG12	1:C:58:VAL:O	2.14	0.48
1:D:173:LYS:O	1:D:176:VAL:HG12	2.13	0.48
1:F:208:TYR:O	1:F:211:SER:HB3	2.13	0.48
1:D:112:LEU:HD12	1:D:136:TRP:HB3	1.94	0.48
1:A:113:SER:HA	1:A:137:ALA:O	2.12	0.48
1:C:177:GLU:O	1:C:180:SER:OG	2.31	0.48
1:E:15:ARG:HG3	1:E:73:SER:HA	1.96	0.48
1:E:2:GLN:OE1	3:E:405:HOH:O	2.20	0.47
1:C:211:SER:OG	1:C:225:ARG:NH1	2.47	0.47
1:F:41:LYS:N	1:F:66:ASP:OD1	2.42	0.47
1:A:193:LYS:O	1:A:222:ASN:ND2	2.45	0.47
1:D:253:LYS:HG2	1:D:257:LEU:HD12	1.96	0.47
1:F:124:ARG:HG3	1:F:190:GLY:H	1.78	0.47
1:A:19:HIS:CE1	1:A:52:SER:HB3	2.49	0.47
1:D:191:TYR:CZ	1:D:193:LYS:HB2	2.49	0.47
1:E:242:LYS:O	1:E:246:GLU:HG3	2.15	0.47
1:E:99:GLN:HB2	1:E:102:THR:OG1	2.15	0.47
1:B:154:TYR:HB3	1:B:158:NLE:HE2	1.97	0.46
1:C:237:SER:O	1:C:241:GLU:HB2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:ILE:HD11	1:D:184:ILE:HD12	1.97	0.46
1:E:87:GLU:OE1	1:E:90:ARG:HD2	2.14	0.46
1:D:201:THR:HG23	1:D:229:GLU:HA	1.97	0.46
1:D:237:SER:O	1:D:241:GLU:HB3	2.16	0.46
1:B:48:PHE:HB3	1:B:64:ARG:HD2	1.98	0.46
1:E:81:THR:HA	1:E:175:PHE:HD1	1.81	0.46
1:F:29:PRO:HB3	1:F:106:ARG:HB3	1.98	0.46
1:E:124:ARG:HD3	1:E:190:GLY:C	2.40	0.46
1:F:35:HIS:CE1	1:F:67:PHE:HD2	2.34	0.46
1:C:37:PHE:O	1:C:38:THR:OG1	2.32	0.46
1:F:83:SER:HA	1:F:86:LEU:HB2	1.97	0.46
1:D:86:LEU:HD22	1:D:122:VAL:HG23	1.98	0.46
1:E:110:LEU:HA	1:E:134:VAL:O	2.16	0.46
1:E:81:THR:H	1:E:84:SER:HB3	1.81	0.45
1:D:191:TYR:CE2	1:D:193:LYS:HB2	2.52	0.45
1:E:1:NLE:N	1:E:56:GLU:OE1	2.48	0.45
1:E:206:VAL:HG22	1:E:210:VAL:HG11	1.97	0.45
1:C:114:NLE:HG2	1:C:142:NLE:HD2	1.98	0.45
1:A:208:TYR:OH	1:A:227:THR:OG1	2.24	0.45
1:E:5:VAL:O	3:E:406:HOH:O	2.21	0.45
1:C:58:VAL:HG11	1:C:253:LYS:HB2	1.99	0.45
1:B:29:PRO:HG3	1:B:257:LEU:CD2	2.43	0.45
1:B:124:ARG:HG3	1:B:189:LYS:O	2.16	0.45
1:C:103:ASP:OD1	1:C:106:ARG:HG3	2.15	0.45
1:A:7:ILE:HB	1:A:14:LEU:HB2	1.98	0.45
1:E:24:VAL:HG12	1:E:28:VAL:HG11	1.98	0.45
1:B:81:THR:HA	1:B:175:PHE:CD2	2.52	0.45
1:E:148:ASN:OD1	1:E:149:GLU:N	2.49	0.45
1:F:127:LYS:O	1:F:193:LYS:NZ	2.48	0.45
1:A:25:LYS:HE3	1:A:25:LYS:HB3	1.72	0.45
1:A:29:PRO:HG3	1:A:106:ARG:HD2	1.98	0.45
1:B:45:HIS:CD2	1:B:236:LYS:HG3	2.52	0.45
1:B:9:TYR:OH	1:B:88:ASP:OD1	2.20	0.44
1:D:230:ASN:O	1:D:240:TRP:CZ2	2.69	0.44
1:E:238:LEU:O	1:E:242:LYS:HB2	2.17	0.44
1:B:105:GLU:O	1:B:131:LYS:HG3	2.17	0.44
1:A:94:LYS:O	1:A:98:GLU:HG3	2.17	0.44
1:B:191:TYR:CZ	1:B:193:LYS:HB2	2.53	0.44
1:E:29:PRO:HB2	1:E:60:ILE:HG23	2.00	0.44
1:B:242:LYS:O	1:B:246:GLU:HB2	2.18	0.44
1:C:175:PHE:O	1:C:179:ILE:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:81:THR:HG23	1:F:84:SER:H	1.83	0.44
1:A:133:LEU:HD23	1:A:195:VAL:HG13	1.99	0.43
1:B:160:GLN:NE2	1:D:99:GLN:OE1	2.47	0.43
1:D:129:GLU:O	3:D:403:HOH:O	2.21	0.43
1:B:137:ALA:HA	1:B:199:HIS:ND1	2.33	0.43
1:D:81:THR:HA	1:D:175:PHE:CD2	2.53	0.43
1:F:210:VAL:O	1:F:214:ILE:HG13	2.19	0.43
1:B:21:PRO:CB	1:B:59:GLY:HA2	2.48	0.43
1:F:3:LYS:HA	1:F:3:LYS:HD3	1.83	0.43
1:E:24:VAL:CG1	1:E:28:VAL:HG11	2.48	0.43
1:F:121:ILE:O	1:F:125:GLU:HG3	2.19	0.43
1:A:81:THR:HA	1:A:175:PHE:CD1	2.51	0.43
1:A:114:NLE:HD2	3:A:407:HOH:O	2.18	0.43
1:C:228:ILE:HG23	1:C:243:LYS:HD2	2.01	0.43
1:D:112:LEU:HA	1:D:136:TRP:O	2.19	0.43
1:A:20:LEU:HD11	1:A:101:THR:HG21	1.99	0.43
1:B:53:ARG:O	1:B:56:GLU:HG2	2.18	0.43
1:B:63:VAL:HG11	1:B:96:VAL:CG2	2.49	0.42
1:D:19:HIS:HB2	1:D:62:SER:OG	2.19	0.42
1:E:173:LYS:HB3	1:E:173:LYS:HE2	1.58	0.42
1:B:196:LEU:HD11	1:B:226:VAL:CG2	2.46	0.42
1:C:6:GLU:OE1	3:C:402:HOH:O	2.22	0.42
1:D:38:THR:HB	1:D:169:HIS:CE1	2.54	0.42
1:D:184:ILE:O	1:D:188:SER:OG	2.25	0.42
1:E:33:NLE:O	3:E:407:HOH:O	2.22	0.42
1:F:112:LEU:HA	1:F:136:TRP:O	2.19	0.42
1:A:64:ARG:HH22	1:B:72:GLU:CD	2.26	0.42
1:B:222:ASN:OD1	1:B:222:ASN:N	2.52	0.42
1:A:40:ASN:OD1	1:A:42:VAL:HG22	2.19	0.42
1:D:222:ASN:OD1	1:D:222:ASN:N	2.51	0.42
1:E:247:GLU:C	3:E:402:HOH:O	2.63	0.42
1:A:23:ASP:OD1	1:A:23:ASP:N	2.40	0.42
1:F:173:LYS:O	1:F:176:VAL:HG22	2.20	0.42
1:A:158:NLE:HB3	1:A:158:NLE:HE3	1.73	0.42
1:D:119:ALA:HB1	1:D:133:LEU:HD11	2.02	0.42
1:E:30:NLE:HD3	1:E:61:GLY:HA3	2.01	0.42
1:C:256:LEU:HD23	1:C:256:LEU:HA	1.83	0.42
1:F:212:ASP:O	1:F:216:LYS:HD3	2.19	0.42
1:F:161:LEU:HD12	1:F:161:LEU:HA	1.82	0.42
1:B:126:TYR:HB3	1:B:129:GLU:HB2	2.02	0.41
1:D:216:LYS:N	1:D:216:LYS:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:243:LYS:HG3	3:D:401:HOH:O	2.20	0.41
1:F:131:LYS:HD3	1:F:255:GLU:O	2.20	0.41
1:C:60:ILE:HD13	1:C:256:LEU:HD12	2.00	0.41
1:C:209:LYS:O	1:C:212:ASP:HB2	2.20	0.41
1:D:32:ILE:HB	1:D:109:LEU:HD23	2.02	0.41
1:B:9:TYR:CD1	1:B:91:GLN:HG3	2.54	0.41
1:D:198:VAL:HG22	1:D:226:VAL:CG2	2.50	0.41
1:A:201:THR:HB	1:A:230:ASN:H	1.85	0.41
1:D:40:ASN:OD1	1:D:42:VAL:HB	2.20	0.41
1:A:99:GLN:HB2	1:A:102:THR:OG1	2.20	0.41
1:A:142:NLE:HB2	1:A:143:PRO:HD3	2.03	0.41
1:B:21:PRO:CG	1:B:28:VAL:HG21	2.47	0.41
1:C:238:LEU:HG	1:D:165:ASP:OD1	2.20	0.41
1:E:51:NLE:HE2	1:E:55:LEU:HD11	2.03	0.41
1:F:86:LEU:HD12	1:F:86:LEU:HA	1.76	0.41
1:B:228:ILE:HG23	1:B:243:LYS:HD3	2.02	0.41
1:D:188:SER:HB2	1:D:218:VAL:HG11	2.02	0.41
1:F:23:ASP:OD1	1:F:23:ASP:N	2.42	0.41
1:F:37:PHE:HZ	1:F:171:LEU:HD13	1.86	0.41
1:F:142:NLE:N	1:F:143:PRO:HD2	2.35	0.41
1:A:121:ILE:HG12	1:A:187:LEU:HB3	2.03	0.41
1:B:121:ILE:HG12	1:B:187:LEU:HB3	2.02	0.41
1:C:9:TYR:OH	1:C:88:ASP:OD1	2.35	0.41
1:C:106:ARG:NH1	1:C:257:LEU:HA	2.35	0.41
1:D:69:GLY:HA2	1:D:75:GLY:C	2.46	0.41
1:C:179:ILE:HG13	1:C:179:ILE:O	2.21	0.41
1:B:86:LEU:HD12	1:B:86:LEU:HA	1.80	0.40
1:E:1:NLE:HE3	1:E:20:LEU:HB2	2.03	0.40
1:A:13:THR:HB	1:A:74:ASP:OD1	2.21	0.40
1:B:147:NLE:HE2	1:B:147:NLE:HB2	1.95	0.40
1:C:165:ASP:CG	1:D:238:LEU:HG	2.46	0.40
1:D:29:PRO:HG3	1:D:256:LEU:O	2.22	0.40
1:A:135:LEU:HB3	1:A:138:PRO:HB3	2.04	0.40
1:B:30:NLE:HG2	1:B:102:THR:HG23	2.04	0.40
1:C:12:LYS:HB3	1:C:68:TYR:CE1	2.57	0.40

All (1) 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:A:90:ARG:NH2	1:C:23:ASP:O[1_545]	2.17	0.03

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	246/267 (92%)	235 (96%)	10 (4%)	1 (0%)	30	60
1	B	245/267 (92%)	233 (95%)	12 (5%)	0	100	100
1	C	235/267 (88%)	226 (96%)	9 (4%)	0	100	100
1	D	234/267 (88%)	226 (97%)	8 (3%)	0	100	100
1	E	237/267 (89%)	227 (96%)	10 (4%)	0	100	100
1	F	239/267 (90%)	227 (95%)	12 (5%)	0	100	100
All	All	1436/1602 (90%)	1374 (96%)	61 (4%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	GLY

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	211/218 (97%)	211 (100%)	0	100	100
1	B	210/218 (96%)	210 (100%)	0	100	100
1	C	202/218 (93%)	202 (100%)	0	100	100
1	D	202/218 (93%)	202 (100%)	0	100	100
1	E	206/218 (94%)	203 (98%)	3 (2%)	57	84
1	F	206/218 (94%)	206 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1237/1308 (95%)	1234 (100%)	3 (0%)	87 96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	180[A]	SER
1	E	180[B]	SER
1	E	228	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	40	ASN
1	D	10	ASN
1	D	169	HIS
1	E	99	GLN
1	E	233	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

63 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	NLE	A	80	1	6,7,8	0.52	0	2,7,9	0.33	0
1	NLE	F	33	1	6,7,8	0.51	0	2,7,9	0.23	0
1	NLE	B	30	1	6,7,8	0.49	0	2,7,9	0.45	0
1	NLE	A	33	1	6,7,8	0.48	0	2,7,9	0.41	0
1	NLE	B	51	1	6,7,8	0.46	0	2,7,9	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	NLE	D	158	1	6,7,8	0.58	0	2,7,9	0.43	0
1	NLE	C	80	1	6,7,8	0.52	0	2,7,9	0.26	0
1	NLE	D	17	1	6,7,8	0.48	0	2,7,9	0.41	0
1	NLE	A	18	1	6,7,8	0.46	0	2,7,9	0.37	0
1	NLE	E	147	1	6,7,8	0.50	0	2,7,9	0.42	0
1	NLE	D	18	1	6,7,8	0.46	0	2,7,9	0.29	0
1	NLE	A	158	1	6,7,8	0.51	0	2,7,9	0.53	0
1	NLE	B	142	1	6,7,8	0.49	0	2,7,9	0.37	0
1	NLE	D	30	1	6,7,8	0.46	0	2,7,9	0.43	0
1	NLE	C	18	1	6,7,8	0.50	0	2,7,9	0.44	0
1	NLE	C	142	1	6,7,8	0.46	0	2,7,9	0.37	0
1	NLE	F	142	1	6,7,8	0.44	0	2,7,9	0.39	0
1	NLE	A	142	1	6,7,8	0.52	0	2,7,9	0.36	0
1	NLE	C	147	1	6,7,8	0.48	0	2,7,9	0.37	0
1	NLE	E	51	1	6,7,8	0.49	0	2,7,9	0.39	0
1	NLE	E	33	1	6,7,8	0.45	0	2,7,9	0.32	0
1	NLE	D	142	1	6,7,8	0.45	0	2,7,9	0.35	0
1	NLE	B	80	1	6,7,8	0.52	0	2,7,9	0.27	0
1	NLE	F	147	1	6,7,8	0.47	0	2,7,9	0.33	0
1	NLE	D	51	1	6,7,8	0.48	0	2,7,9	0.36	0
1	NLE	F	1	1	6,7,8	0.50	0	2,7,9	0.36	0
1	NLE	E	17	1	6,7,8	0.46	0	2,7,9	0.40	0
1	NLE	C	1	1	6,7,8	0.42	0	2,7,9	0.42	0
1	NLE	F	17	1	6,7,8	0.49	0	2,7,9	0.40	0
1	NLE	A	1	1	6,7,8	0.49	0	2,7,9	0.39	0
1	NLE	F	114	1	6,7,8	0.50	0	2,7,9	0.42	0
1	NLE	E	80	1	6,7,8	0.52	0	2,7,9	0.32	0
1	NLE	B	147	1	6,7,8	0.49	0	2,7,9	0.29	0
1	NLE	E	114	1	6,7,8	0.46	0	2,7,9	0.33	0
1	NLE	B	18	1	6,7,8	0.48	0	2,7,9	0.33	0
1	NLE	A	147	1	6,7,8	0.48	0	2,7,9	0.41	0
1	NLE	D	33	1	6,7,8	0.49	0	2,7,9	0.41	0
1	NLE	B	114	1	6,7,8	0.49	0	2,7,9	0.37	0
1	NLE	E	142	1	6,7,8	0.47	0	2,7,9	0.37	0
1	NLE	F	158	1	6,7,8	0.67	0	2,7,9	0.70	0
1	NLE	C	17	1	6,7,8	0.49	0	2,7,9	0.44	0
1	NLE	C	114	1	6,7,8	0.49	0	2,7,9	0.35	0
1	NLE	F	18	1	6,7,8	0.47	0	2,7,9	0.27	0
1	NLE	E	30	1	6,7,8	0.45	0	2,7,9	0.43	0
1	NLE	C	33	1	6,7,8	0.48	0	2,7,9	0.42	0
1	NLE	D	114	1	6,7,8	0.48	0	2,7,9	0.43	0
1	NLE	B	158	1	6,7,8	0.46	0	2,7,9	0.49	0
1	NLE	F	30	1	6,7,8	0.49	0	2,7,9	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	NLE	C	30	1	6,7,8	0.48	0	2,7,9	0.41	0
1	NLE	A	30	1	6,7,8	0.43	0	2,7,9	0.46	0
1	NLE	A	17	1	6,7,8	0.47	0	2,7,9	0.42	0
1	NLE	B	1	1	6,7,8	0.48	0	2,7,9	0.41	0
1	NLE	A	114	1	6,7,8	0.49	0	2,7,9	0.33	0
1	NLE	F	51	1	6,7,8	0.50	0	2,7,9	0.40	0
1	NLE	B	33	1	6,7,8	0.45	0	2,7,9	0.44	0
1	NLE	D	80	1	6,7,8	0.52	0	2,7,9	0.28	0
1	NLE	E	1	1	6,7,8	0.48	0	2,7,9	0.39	0
1	NLE	A	51	1	6,7,8	0.49	0	2,7,9	0.35	0
1	NLE	E	18	1	6,7,8	0.46	0	2,7,9	0.43	0
1	NLE	D	1	1	6,7,8	0.49	0	2,7,9	0.48	0
1	NLE	C	51	1	6,7,8	0.51	0	2,7,9	0.39	0
1	NLE	F	80	1	6,7,8	0.50	0	2,7,9	0.31	0
1	NLE	B	17	1	6,7,8	0.46	0	2,7,9	0.41	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
1	NLE	A	80	1	-	2/5/6/8	-
1	NLE	F	33	1	-	0/5/6/8	-
1	NLE	B	30	1	-	1/5/6/8	-
1	NLE	A	33	1	-	0/5/6/8	-
1	NLE	B	51	1	-	0/5/6/8	-
1	NLE	D	158	1	-	1/5/6/8	-
1	NLE	C	80	1	-	1/5/6/8	-
1	NLE	D	17	1	-	2/5/6/8	-
1	NLE	A	18	1	-	0/5/6/8	-
1	NLE	E	147	1	-	1/5/6/8	-
1	NLE	D	18	1	-	1/5/6/8	-
1	NLE	A	158	1	-	4/5/6/8	-
1	NLE	B	142	1	-	1/5/6/8	-
1	NLE	D	30	1	-	1/5/6/8	-
1	NLE	C	18	1	-	0/5/6/8	-
1	NLE	C	142	1	-	1/5/6/8	-
1	NLE	F	142	1	-	1/5/6/8	-
1	NLE	A	142	1	-	1/5/6/8	-
1	NLE	C	147	1	-	1/5/6/8	-
1	NLE	E	51	1	-	0/5/6/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	NLE	E	33	1	-	0/5/6/8	-
1	NLE	D	142	1	-	1/5/6/8	-
1	NLE	B	80	1	-	1/5/6/8	-
1	NLE	F	147	1	-	2/5/6/8	-
1	NLE	D	51	1	-	2/5/6/8	-
1	NLE	F	1	1	-	1/5/6/8	-
1	NLE	E	17	1	-	2/5/6/8	-
1	NLE	C	1	1	-	1/5/6/8	-
1	NLE	F	17	1	-	2/5/6/8	-
1	NLE	A	1	1	-	1/5/6/8	-
1	NLE	F	114	1	-	0/5/6/8	-
1	NLE	E	80	1	-	1/5/6/8	-
1	NLE	B	147	1	-	1/5/6/8	-
1	NLE	E	114	1	-	0/5/6/8	-
1	NLE	B	18	1	-	0/5/6/8	-
1	NLE	A	147	1	-	3/5/6/8	-
1	NLE	D	33	1	-	0/5/6/8	-
1	NLE	B	114	1	-	1/5/6/8	-
1	NLE	E	142	1	-	0/5/6/8	-
1	NLE	F	158	1	-	0/5/6/8	-
1	NLE	C	17	1	-	4/5/6/8	-
1	NLE	C	114	1	-	1/5/6/8	-
1	NLE	F	18	1	-	1/5/6/8	-
1	NLE	E	30	1	-	1/5/6/8	-
1	NLE	C	33	1	-	0/5/6/8	-
1	NLE	D	114	1	-	1/5/6/8	-
1	NLE	B	158	1	-	2/5/6/8	-
1	NLE	F	30	1	-	0/5/6/8	-
1	NLE	C	30	1	-	2/5/6/8	-
1	NLE	A	30	1	-	2/5/6/8	-
1	NLE	A	17	1	-	3/5/6/8	-
1	NLE	B	1	1	-	1/5/6/8	-
1	NLE	A	114	1	-	0/5/6/8	-
1	NLE	F	51	1	-	0/5/6/8	-
1	NLE	B	33	1	-	0/5/6/8	-
1	NLE	D	80	1	-	2/5/6/8	-
1	NLE	E	1	1	-	1/5/6/8	-
1	NLE	A	51	1	-	0/5/6/8	-
1	NLE	E	18	1	-	0/5/6/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	NLE	D	1	1	-	2/5/6/8	-
1	NLE	C	51	1	-	0/5/6/8	-
1	NLE	F	80	1	-	2/5/6/8	-
1	NLE	B	17	1	-	0/5/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	F	1	NLE	O-C-CA-CB
1	A	17	NLE	N-CA-CB-CG
1	A	17	NLE	C-CA-CB-CG
1	C	17	NLE	N-CA-CB-CG
1	C	17	NLE	C-CA-CB-CG
1	D	17	NLE	N-CA-CB-CG
1	D	17	NLE	C-CA-CB-CG
1	E	17	NLE	N-CA-CB-CG
1	E	17	NLE	C-CA-CB-CG
1	F	17	NLE	N-CA-CB-CG
1	F	17	NLE	C-CA-CB-CG
1	C	30	NLE	N-CA-CB-CG
1	C	30	NLE	C-CA-CB-CG
1	D	51	NLE	N-CA-CB-CG
1	D	51	NLE	C-CA-CB-CG
1	A	80	NLE	O-C-CA-CB
1	B	114	NLE	O-C-CA-CB
1	C	114	NLE	O-C-CA-CB
1	D	114	NLE	O-C-CA-CB
1	F	147	NLE	C-CA-CB-CG
1	A	1	NLE	CA-CB-CG-CD
1	A	147	NLE	CA-CB-CG-CD
1	C	147	NLE	CE-CD-CG-CB
1	D	158	NLE	CE-CD-CG-CB
1	C	142	NLE	CA-CB-CG-CD
1	C	17	NLE	CA-CB-CG-CD
1	B	142	NLE	CA-CB-CG-CD
1	D	80	NLE	CE-CD-CG-CB
1	A	80	NLE	CA-CB-CG-CD

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Mol	Chain	Res	Type	Atoms
1	C	80	NLE	CA-CB-CG-CD
1	D	80	NLE	CA-CB-CG-CD
1	E	80	NLE	CA-CB-CG-CD
1	F	80	NLE	CA-CB-CG-CD
1	A	158	NLE	CE-CD-CG-CB
1	B	1	NLE	CE-CD-CG-CB
1	E	147	NLE	CE-CD-CG-CB
1	A	30	NLE	CA-CB-CG-CD
1	D	1	NLE	CA-CB-CG-CD
1	A	158	NLE	CA-CB-CG-CD
1	D	18	NLE	CA-CB-CG-CD
1	E	1	NLE	CA-CB-CG-CD
1	A	158	NLE	C-CA-CB-CG
1	B	80	NLE	CA-CB-CG-CD
1	A	147	NLE	N-CA-CB-CG
1	A	158	NLE	N-CA-CB-CG
1	D	142	NLE	CA-CB-CG-CD
1	B	147	NLE	CA-CB-CG-CD
1	A	17	NLE	CA-CB-CG-CD
1	A	142	NLE	CA-CB-CG-CD
1	E	30	NLE	CE-CD-CG-CB
1	F	147	NLE	CE-CD-CG-CB
1	F	142	NLE	CA-CB-CG-CD
1	B	30	NLE	CE-CD-CG-CB
1	C	17	NLE	CE-CD-CG-CB
1	F	80	NLE	CE-CD-CG-CB
1	D	30	NLE	CE-CD-CG-CB
1	F	18	NLE	CE-CD-CG-CB
1	A	30	NLE	C-CA-CB-CG
1	B	158	NLE	CA-CB-CG-CD
1	A	147	NLE	CE-CD-CG-CB
1	B	158	NLE	CE-CD-CG-CB
1	C	1	NLE	N-CA-CB-CG
1	D	1	NLE	CE-CD-CG-CB

There are no ring outliers.

26 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	F	33	NLE	1	0
1	B	30	NLE	1	0
1	B	51	NLE	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	158	NLE	1	0
1	B	142	NLE	1	0
1	D	30	NLE	1	0
1	C	142	NLE	1	0
1	F	142	NLE	1	0
1	A	142	NLE	1	0
1	E	51	NLE	1	0
1	E	33	NLE	1	0
1	C	1	NLE	1	0
1	B	147	NLE	1	0
1	E	114	NLE	1	0
1	B	114	NLE	1	0
1	F	158	NLE	2	0
1	C	17	NLE	1	0
1	C	114	NLE	2	0
1	E	30	NLE	1	0
1	C	33	NLE	1	0
1	B	158	NLE	1	0
1	C	30	NLE	1	0
1	A	17	NLE	1	0
1	A	114	NLE	2	0
1	E	1	NLE	2	0
1	D	1	NLE	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	301	-	4,4,4	0.23	0	6,6,6	0.12	0
2	SO4	C	301	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	E	301	-	4,4,4	0.22	0	6,6,6	0.23	0
2	SO4	F	301	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	B	301	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	D	301	-	4,4,4	0.24	0	6,6,6	0.07	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.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	246/267 (92%)	-0.12	1 (0%) 88 84	24, 45, 68, 95	1 (0%)
1	B	246/267 (92%)	-0.08	1 (0%) 88 84	32, 52, 81, 96	0
1	C	238/267 (89%)	-0.16	1 (0%) 88 84	23, 44, 66, 102	0
1	D	237/267 (88%)	0.17	5 (2%) 63 54	23, 50, 75, 87	0
1	E	239/267 (89%)	-0.19	1 (0%) 88 84	26, 46, 67, 104	1 (0%)
1	F	242/267 (90%)	-0.03	3 (1%) 76 68	31, 50, 78, 98	0
All	All	1448/1602 (90%)	-0.07	12 (0%) 82 75	23, 48, 75, 104	2 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	239	GLU	3.3
1	F	162	GLY	2.8
1	A	108	GLY	2.7
1	E	162	GLY	2.6
1	C	161	LEU	2.6
1	D	157	ILE	2.5
1	D	156	ALA	2.4
1	D	238	LEU	2.2
1	F	146	ILE	2.1
1	B	129	GLU	2.0
1	F	143	PRO	2.0
1	D	143	PRO	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	NLE	C	147	8/9	0.69	0.13	60,74,79,79	0
1	NLE	B	1	8/9	0.76	0.16	62,72,77,79	0
1	NLE	F	158	8/9	0.79	0.20	72,84,93,94	0
1	NLE	D	158	8/9	0.80	0.12	71,77,86,90	0
1	NLE	C	1	8/9	0.84	0.12	37,42,54,57	0
1	NLE	E	147	8/9	0.84	0.10	68,74,89,92	0
1	NLE	D	142	8/9	0.85	0.11	62,64,68,72	0
1	NLE	A	147	8/9	0.86	0.14	53,57,61,71	0
1	NLE	A	18	8/9	0.87	0.12	39,46,51,52	0
1	NLE	B	30	8/9	0.88	0.09	57,61,68,69	0
1	NLE	F	147	8/9	0.88	0.12	73,80,89,90	0
1	NLE	E	1	8/9	0.90	0.12	41,45,52,56	0
1	NLE	D	80	8/9	0.90	0.08	45,46,56,58	0
1	NLE	A	1	8/9	0.90	0.15	54,65,72,74	0
1	NLE	F	80	8/9	0.91	0.10	39,45,52,56	0
1	NLE	F	142	8/9	0.91	0.11	56,59,67,70	0
1	NLE	B	158	8/9	0.91	0.13	48,50,58,59	0
1	NLE	D	114	8/9	0.91	0.08	37,44,45,47	0
1	NLE	F	114	8/9	0.91	0.10	37,42,47,47	0
1	NLE	D	18	8/9	0.92	0.10	20,26,32,33	0
1	NLE	E	51	8/9	0.92	0.10	28,35,37,37	0
1	NLE	C	80	8/9	0.92	0.08	36,39,49,52	0
1	NLE	A	30	8/9	0.92	0.13	40,46,49,53	0
1	NLE	A	114	8/9	0.93	0.11	29,34,35,39	0
1	NLE	A	158	8/9	0.93	0.10	58,60,61,66	0
1	NLE	A	17	8/9	0.93	0.11	40,44,46,47	0
1	NLE	B	18	8/9	0.93	0.09	50,54,57,60	0
1	NLE	B	17	8/9	0.93	0.12	46,49,50,51	0
1	NLE	C	17	8/9	0.94	0.10	23,26,31,37	0
1	NLE	E	80	8/9	0.94	0.07	41,43,48,48	0
1	NLE	B	147	8/9	0.94	0.08	48,52,53,63	0
1	NLE	E	33	8/9	0.94	0.11	23,33,36,41	0
1	NLE	E	18	8/9	0.94	0.09	29,37,41,49	0
1	NLE	B	114	8/9	0.94	0.08	35,39,43,44	0
1	NLE	C	114	8/9	0.94	0.08	25,27,36,39	0
1	NLE	A	80	8/9	0.94	0.09	34,39,43,44	0
1	NLE	B	80	8/9	0.94	0.08	34,40,46,48	0
1	NLE	E	17	8/9	0.94	0.10	29,34,36,39	0
1	NLE	A	33	8/9	0.95	0.08	29,35,40,41	0
1	NLE	B	33	8/9	0.95	0.08	43,49,50,54	0
1	NLE	D	1	8/9	0.95	0.09	34,37,41,43	0
1	NLE	F	33	8/9	0.95	0.10	34,37,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	NLE	E	114	8/9	0.95	0.07	29,33,38,38	0
1	NLE	C	51	8/9	0.95	0.07	27,35,36,36	0
1	NLE	A	142	8/9	0.95	0.09	33,37,50,50	0
1	NLE	B	142	8/9	0.95	0.08	30,37,42,42	0
1	NLE	F	1	8/9	0.95	0.10	39,45,48,51	0
1	NLE	E	142	8/9	0.95	0.14	46,51,54,55	0
1	NLE	D	51	8/9	0.96	0.07	35,37,43,43	0
1	NLE	E	30	8/9	0.96	0.07	43,46,49,52	0
1	NLE	F	51	8/9	0.96	0.07	36,39,43,43	0
1	NLE	C	142	8/9	0.96	0.14	49,52,56,56	0
1	NLE	F	18	8/9	0.96	0.07	29,33,36,42	0
1	NLE	A	51	8/9	0.96	0.09	33,37,41,43	0
1	NLE	B	51	8/9	0.96	0.10	45,50,53,53	0
1	NLE	C	30	8/9	0.96	0.08	40,43,46,46	0
1	NLE	D	33	8/9	0.97	0.09	28,36,41,43	0
1	NLE	D	30	8/9	0.97	0.08	27,34,39,43	0
1	NLE	F	30	8/9	0.97	0.07	39,41,44,44	0
1	NLE	C	33	8/9	0.97	0.10	27,32,33,34	0
1	NLE	F	17	8/9	0.98	0.07	27,32,34,41	0
1	NLE	C	18	8/9	0.98	0.06	26,30,32,33	0
1	NLE	D	17	8/9	0.99	0.04	19,25,26,29	0

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(\AA^2)	Q<0.9
2	SO4	E	301	5/5	0.86	0.11	68,78,83,88	0
2	SO4	B	301	5/5	0.90	0.07	55,55,65,73	0
2	SO4	D	301	5/5	0.95	0.10	54,56,73,80	0
2	SO4	C	301	5/5	0.95	0.10	63,69,73,84	0
2	SO4	F	301	5/5	0.95	0.07	51,59,65,77	0
2	SO4	A	301	5/5	0.96	0.07	43,60,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.