



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2021 – 11:58 PM EDT

PDB ID : 3LGV
Title : H198P mutant of the DegS-deltaPDZ protease
Authors : Sohn, J.; Grant, R.A.; Sauer, R.T.
Deposited on : 2010-01-21
Resolution : 2.73 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

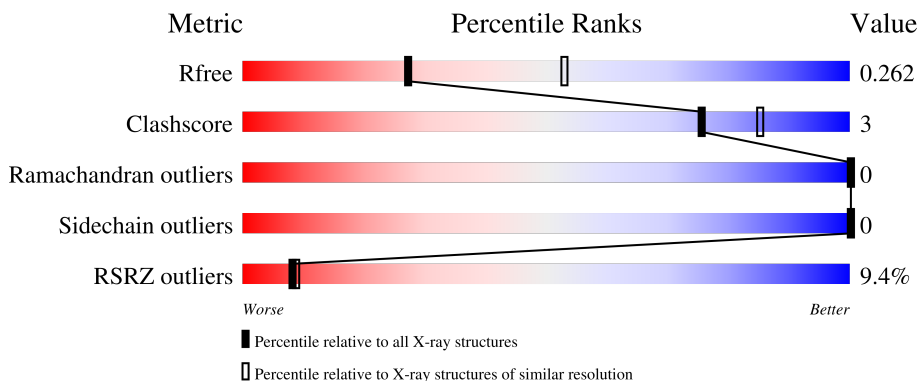
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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.



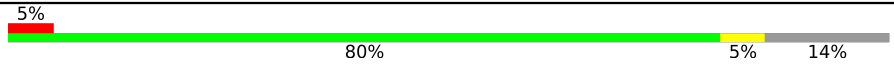
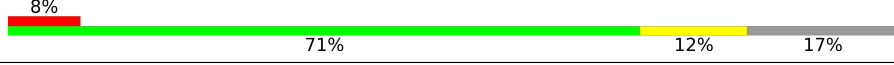
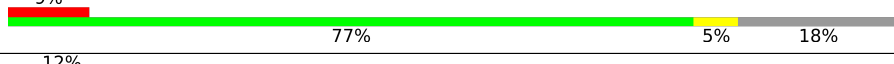

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1271 (2.76-2.72)
Clashscore	141614	1322 (2.76-2.72)
Ramachandran outliers	138981	1297 (2.76-2.72)
Sidechain outliers	138945	1298 (2.76-2.72)
RSRZ outliers	127900	1243 (2.76-2.72)

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	241	 6% 80% 5% 15%
1	B	241	 9% 79% 7% 15%
1	C	241	 7% 79% 6% 15%
1	D	241	 9% 78% 6% 16%
1	E	241	 6% 73% 8% 19%

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Mol	Chain	Length	Quality of chain
1	F	241	
1	G	241	
1	H	241	
1	I	241	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 27099 atoms, of which 13516 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 Protease degS.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	204	3026	946	1518	263	296	3	0	0	0
1	B	206	3079	957	1550	272	297	3	0	0	0
1	C	205	3064	951	1542	272	296	3	0	0	0
1	D	203	3026	942	1523	268	290	3	0	0	0
1	E	195	2910	912	1463	253	279	3	0	0	0
1	F	207	3069	957	1541	272	296	3	0	0	0
1	G	200	2962	925	1490	258	286	3	0	0	0
1	H	197	2902	907	1461	254	277	3	0	0	0
1	I	193	2853	893	1428	250	279	3	0	0	0

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	expression tag	UNP P0AEE3
A	17	ARG	-	expression tag	UNP P0AEE3
A	18	GLY	-	expression tag	UNP P0AEE3
A	19	SER	-	expression tag	UNP P0AEE3
A	20	HIS	-	expression tag	UNP P0AEE3
A	21	HIS	-	expression tag	UNP P0AEE3
A	22	HIS	-	expression tag	UNP P0AEE3
A	23	HIS	-	expression tag	UNP P0AEE3
A	24	HIS	-	expression tag	UNP P0AEE3
A	25	HIS	-	expression tag	UNP P0AEE3
A	26	GLY	-	expression tag	UNP P0AEE3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	198	PRO	HIS	engineered mutation	UNP P0AEE3
B	16	MET	-	expression tag	UNP P0AEE3
B	17	ARG	-	expression tag	UNP P0AEE3
B	18	GLY	-	expression tag	UNP P0AEE3
B	19	SER	-	expression tag	UNP P0AEE3
B	20	HIS	-	expression tag	UNP P0AEE3
B	21	HIS	-	expression tag	UNP P0AEE3
B	22	HIS	-	expression tag	UNP P0AEE3
B	23	HIS	-	expression tag	UNP P0AEE3
B	24	HIS	-	expression tag	UNP P0AEE3
B	25	HIS	-	expression tag	UNP P0AEE3
B	26	GLY	-	expression tag	UNP P0AEE3
B	198	PRO	HIS	engineered mutation	UNP P0AEE3
C	16	MET	-	expression tag	UNP P0AEE3
C	17	ARG	-	expression tag	UNP P0AEE3
C	18	GLY	-	expression tag	UNP P0AEE3
C	19	SER	-	expression tag	UNP P0AEE3
C	20	HIS	-	expression tag	UNP P0AEE3
C	21	HIS	-	expression tag	UNP P0AEE3
C	22	HIS	-	expression tag	UNP P0AEE3
C	23	HIS	-	expression tag	UNP P0AEE3
C	24	HIS	-	expression tag	UNP P0AEE3
C	25	HIS	-	expression tag	UNP P0AEE3
C	26	GLY	-	expression tag	UNP P0AEE3
C	198	PRO	HIS	engineered mutation	UNP P0AEE3
D	16	MET	-	expression tag	UNP P0AEE3
D	17	ARG	-	expression tag	UNP P0AEE3
D	18	GLY	-	expression tag	UNP P0AEE3
D	19	SER	-	expression tag	UNP P0AEE3
D	20	HIS	-	expression tag	UNP P0AEE3
D	21	HIS	-	expression tag	UNP P0AEE3
D	22	HIS	-	expression tag	UNP P0AEE3
D	23	HIS	-	expression tag	UNP P0AEE3
D	24	HIS	-	expression tag	UNP P0AEE3
D	25	HIS	-	expression tag	UNP P0AEE3
D	26	GLY	-	expression tag	UNP P0AEE3
D	198	PRO	HIS	engineered mutation	UNP P0AEE3
E	16	MET	-	expression tag	UNP P0AEE3
E	17	ARG	-	expression tag	UNP P0AEE3
E	18	GLY	-	expression tag	UNP P0AEE3
E	19	SER	-	expression tag	UNP P0AEE3
E	20	HIS	-	expression tag	UNP P0AEE3

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Chain	Residue	Modelled	Actual	Comment	Reference
E	21	HIS	-	expression tag	UNP P0AEE3
E	22	HIS	-	expression tag	UNP P0AEE3
E	23	HIS	-	expression tag	UNP P0AEE3
E	24	HIS	-	expression tag	UNP P0AEE3
E	25	HIS	-	expression tag	UNP P0AEE3
E	26	GLY	-	expression tag	UNP P0AEE3
E	198	PRO	HIS	engineered mutation	UNP P0AEE3
F	16	MET	-	expression tag	UNP P0AEE3
F	17	ARG	-	expression tag	UNP P0AEE3
F	18	GLY	-	expression tag	UNP P0AEE3
F	19	SER	-	expression tag	UNP P0AEE3
F	20	HIS	-	expression tag	UNP P0AEE3
F	21	HIS	-	expression tag	UNP P0AEE3
F	22	HIS	-	expression tag	UNP P0AEE3
F	23	HIS	-	expression tag	UNP P0AEE3
F	24	HIS	-	expression tag	UNP P0AEE3
F	25	HIS	-	expression tag	UNP P0AEE3
F	26	GLY	-	expression tag	UNP P0AEE3
F	198	PRO	HIS	engineered mutation	UNP P0AEE3
G	16	MET	-	expression tag	UNP P0AEE3
G	17	ARG	-	expression tag	UNP P0AEE3
G	18	GLY	-	expression tag	UNP P0AEE3
G	19	SER	-	expression tag	UNP P0AEE3
G	20	HIS	-	expression tag	UNP P0AEE3
G	21	HIS	-	expression tag	UNP P0AEE3
G	22	HIS	-	expression tag	UNP P0AEE3
G	23	HIS	-	expression tag	UNP P0AEE3
G	24	HIS	-	expression tag	UNP P0AEE3
G	25	HIS	-	expression tag	UNP P0AEE3
G	26	GLY	-	expression tag	UNP P0AEE3
G	198	PRO	HIS	engineered mutation	UNP P0AEE3
H	16	MET	-	expression tag	UNP P0AEE3
H	17	ARG	-	expression tag	UNP P0AEE3
H	18	GLY	-	expression tag	UNP P0AEE3
H	19	SER	-	expression tag	UNP P0AEE3
H	20	HIS	-	expression tag	UNP P0AEE3
H	21	HIS	-	expression tag	UNP P0AEE3
H	22	HIS	-	expression tag	UNP P0AEE3
H	23	HIS	-	expression tag	UNP P0AEE3
H	24	HIS	-	expression tag	UNP P0AEE3
H	25	HIS	-	expression tag	UNP P0AEE3
H	26	GLY	-	expression tag	UNP P0AEE3

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Chain	Residue	Modelled	Actual	Comment	Reference
H	198	PRO	HIS	engineered mutation	UNP P0AEE3
I	16	MET	-	expression tag	UNP P0AEE3
I	17	ARG	-	expression tag	UNP P0AEE3
I	18	GLY	-	expression tag	UNP P0AEE3
I	19	SER	-	expression tag	UNP P0AEE3
I	20	HIS	-	expression tag	UNP P0AEE3
I	21	HIS	-	expression tag	UNP P0AEE3
I	22	HIS	-	expression tag	UNP P0AEE3
I	23	HIS	-	expression tag	UNP P0AEE3
I	24	HIS	-	expression tag	UNP P0AEE3
I	25	HIS	-	expression tag	UNP P0AEE3
I	26	GLY	-	expression tag	UNP P0AEE3
I	198	PRO	HIS	engineered mutation	UNP P0AEE3

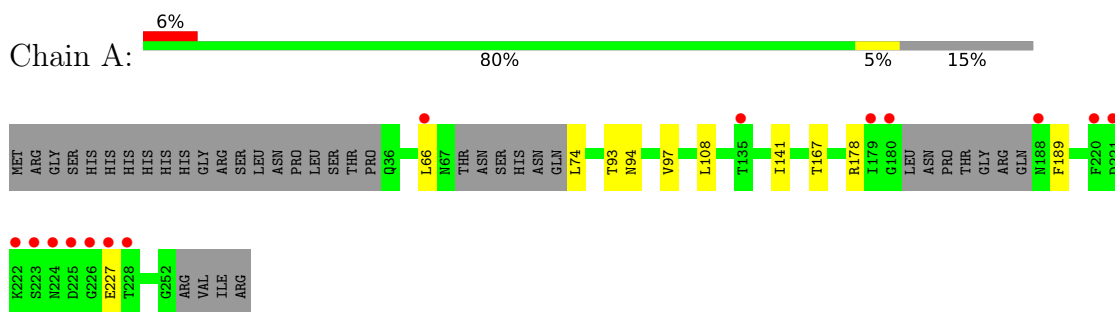
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	41	Total O 41 41	0	0
2	B	30	Total O 30 30	0	0
2	C	28	Total O 28 28	0	0
2	D	17	Total O 17 17	0	0
2	E	21	Total O 21 21	0	0
2	F	28	Total O 28 28	0	0
2	G	20	Total O 20 20	0	0
2	H	13	Total O 13 13	0	0
2	I	10	Total O 10 10	0	0

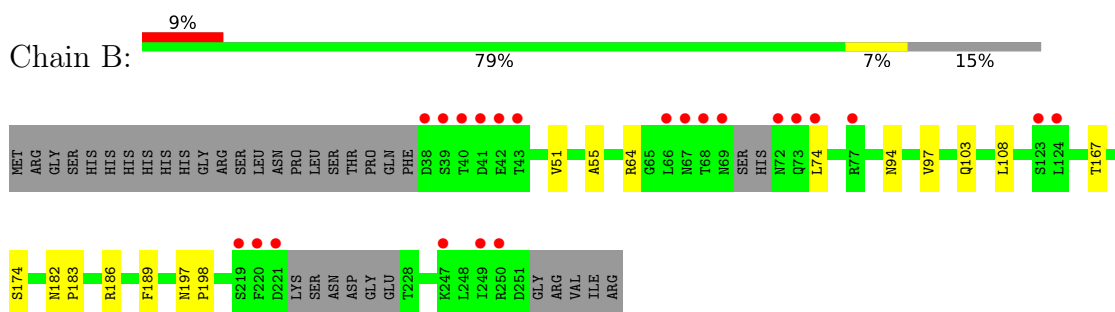
3 Residue-property plots [i](#)

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.

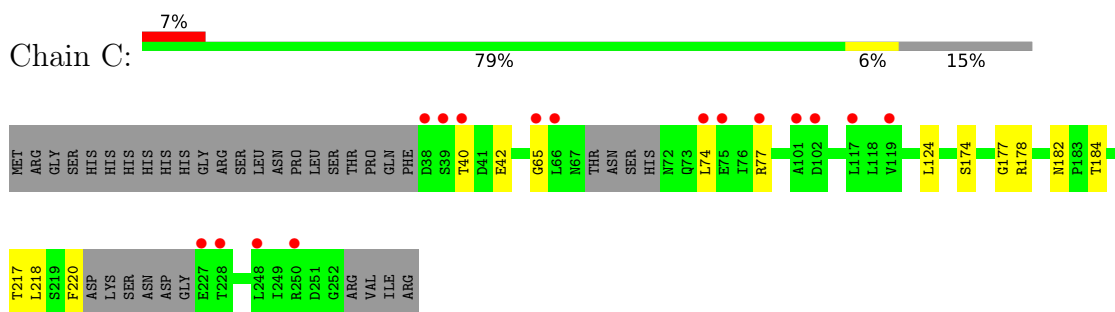
- Molecule 1: Protease degS



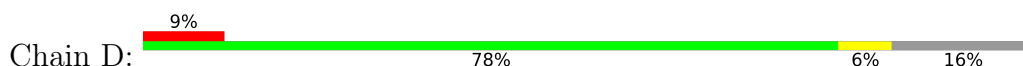
- Molecule 1: Protease degS

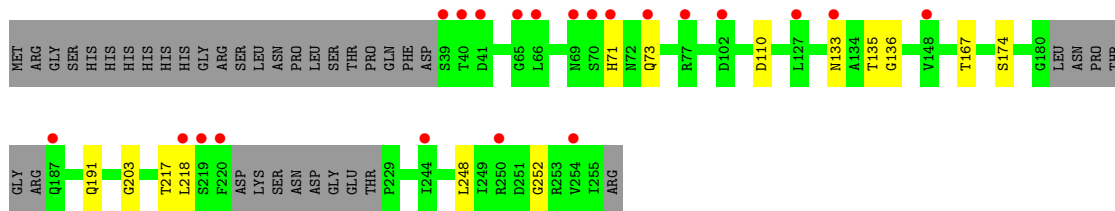


- Molecule 1: Protease degS

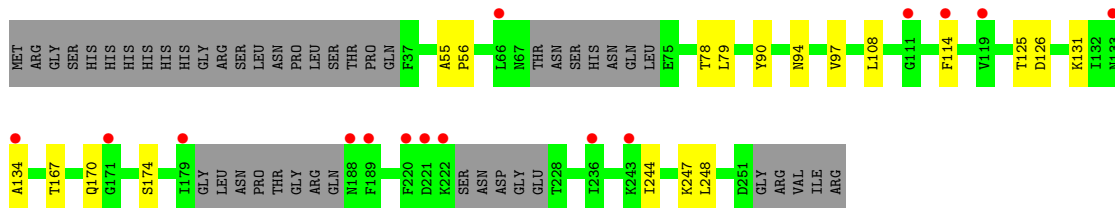
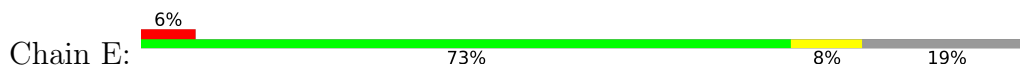


- Molecule 1: Protease degS

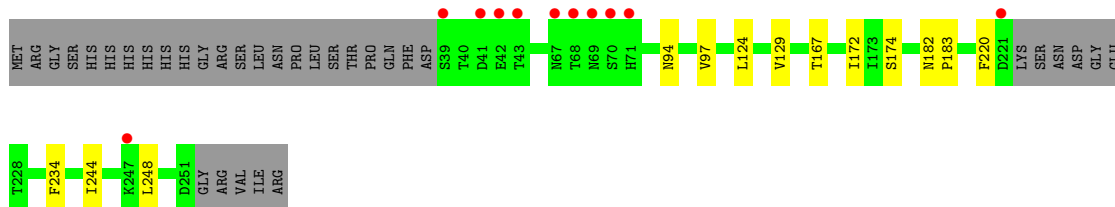
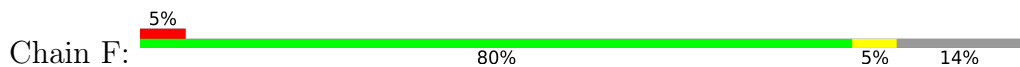




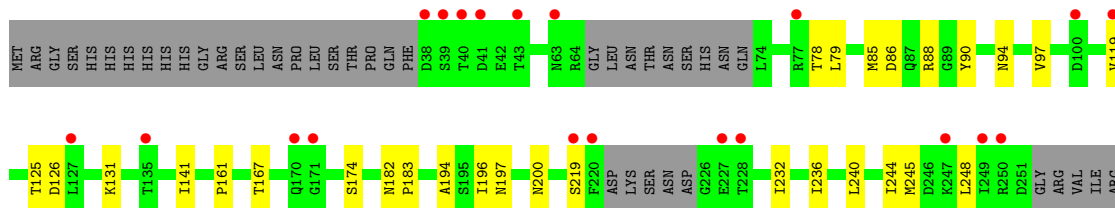
• Molecule 1: Protease degS



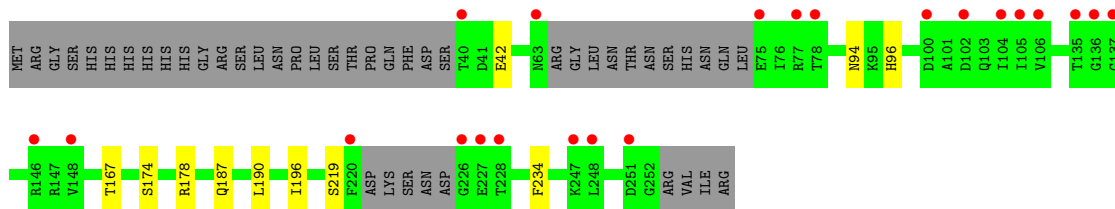
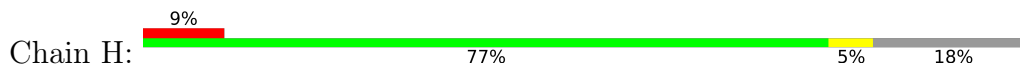
• Molecule 1: Protease degS



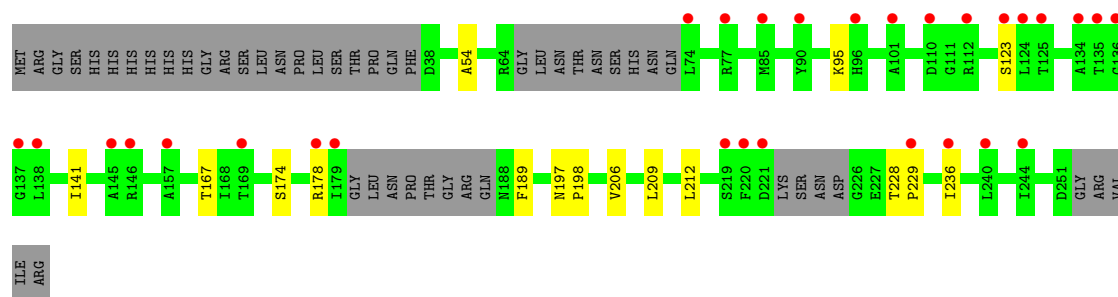
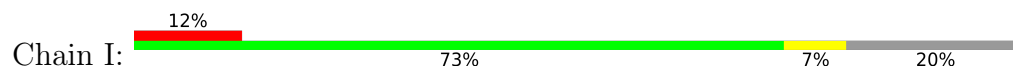
• Molecule 1: Protease degS



• Molecule 1: Protease degS



- Molecule 1: Protease degS



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.54Å 133.56Å 230.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.82 – 2.73 33.82 – 2.73	Depositor EDS
% Data completeness (in resolution range)	94.1 (33.82-2.73) 94.2 (33.82-2.73)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.72Å)	Xtrriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.219 , 0.270 0.210 , 0.262	Depositor DCC
R_{free} test set	2856 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtrriage
Anisotropy	0.124	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	27099	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.15 % 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 1.8376e-04. 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 [i](#)

5.1 Standard geometry [i](#)

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	0.28	0/1526	0.45	0/2073
1	B	0.28	0/1547	0.46	0/2104
1	C	0.29	0/1540	0.46	0/2093
1	D	0.27	0/1521	0.44	0/2068
1	E	0.27	0/1464	0.43	0/1988
1	F	0.28	0/1548	0.44	0/2108
1	G	0.27	0/1490	0.46	0/2028
1	H	0.26	0/1459	0.44	0/1986
1	I	0.26	0/1441	0.46	0/1959
All	All	0.27	0/13536	0.45	0/18407

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	1508	1518	1518	8	0
1	B	1529	1550	1550	10	0
1	C	1522	1542	1541	10	0
1	D	1503	1523	1523	11	0
1	E	1447	1463	1462	13	0
1	F	1528	1541	1541	9	0
1	G	1472	1490	1490	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	1441	1461	1461	8	0
1	I	1425	1428	1428	10	0
2	A	41	0	0	0	0
2	B	30	0	0	0	0
2	C	28	0	0	1	0
2	D	17	0	0	2	0
2	E	21	0	0	0	0
2	F	28	0	0	0	0
2	G	20	0	0	0	0
2	H	13	0	0	0	0
2	I	10	0	0	0	0
All	All	13583	13516	13514	83	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 3.

All (83) 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:G:236:ILE:HG23	1:G:240:LEU:HD23	1.92	0.52
1:H:167:THR:HG23	1:I:174:SER:HB3	1.92	0.51
1:G:161:PRO:HB3	1:G:197:ASN:HB2	1.92	0.50
1:I:206:VAL:HG12	1:I:212:LEU:HA	1.95	0.48
1:B:94:ASN:HB3	1:B:97:VAL:HG23	1.96	0.48
1:B:167:THR:HG23	1:C:174:SER:HB3	1.94	0.48
1:I:197:ASN:HB3	1:I:198:PRO:HD2	1.96	0.47
1:D:174:SER:HB2	1:D:191:GLN:HG2	1.98	0.46
1:D:174:SER:HB3	1:F:167:THR:HG23	1.95	0.46
1:E:167:THR:HG23	1:F:174:SER:HB3	1.97	0.46
1:G:174:SER:HB3	1:I:167:THR:HG23	1.97	0.46
1:D:71:HIS:CD2	1:D:73:GLN:HB2	2.51	0.46
1:C:124:LEU:HD12	1:D:110:ASP:O	2.16	0.45
1:C:65:GLY:HA3	1:C:77:ARG:HD2	1.98	0.45
1:E:247:LYS:HG3	1:E:248:LEU:N	2.32	0.45
1:I:95:LYS:HE3	1:I:123:SER:HA	1.98	0.45
1:A:66:LEU:HD23	1:A:74:LEU:HD23	1.99	0.44
1:B:108:LEU:HD12	1:B:108:LEU:N	2.32	0.44
1:E:125:THR:O	1:E:126:ASP:HB3	2.17	0.44
1:F:94:ASN:HB3	1:F:97:VAL:HG23	1.99	0.44
1:F:124:LEU:HD23	1:F:124:LEU:O	2.18	0.44
1:G:94:ASN:HB3	1:G:97:VAL:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:THR:HG23	1:A:94:ASN:N	2.33	0.44
1:C:217:THR:HG22	1:C:218:LEU:N	2.33	0.44
1:A:178:ARG:O	1:A:189:PHE:HB2	2.18	0.44
1:E:94:ASN:HB3	1:E:97:VAL:HG23	2.00	0.44
1:G:219:SER:HB2	1:G:232:ILE:O	2.17	0.44
1:A:167:THR:HG23	1:B:174:SER:HB3	2.00	0.44
1:I:178:ARG:HA	1:I:178:ARG:HD3	1.79	0.44
1:B:186:ARG:HD2	1:B:189:PHE:CE2	2.53	0.43
1:D:135:THR:HG22	1:D:136:GLY:N	2.32	0.43
1:E:134:ALA:HA	1:H:42:GLU:OE1	2.18	0.43
1:I:228:THR:N	1:I:229:PRO:HD3	2.32	0.43
1:F:220:PHE:HB2	1:F:234:PHE:HE2	1.82	0.43
1:G:119:VAL:HG12	1:G:119:VAL:O	2.18	0.43
1:A:94:ASN:HB3	1:A:97:VAL:HG23	2.00	0.43
1:A:108:LEU:HD12	1:A:108:LEU:N	2.34	0.43
1:D:248:LEU:O	1:D:252:GLY:HA2	2.18	0.43
1:G:244:ILE:O	1:G:248:LEU:HG	2.18	0.43
1:G:161:PRO:HG3	1:G:200:ASN:OD1	2.19	0.43
1:D:217:THR:O	1:D:218:LEU:HG	2.18	0.43
1:E:78:THR:HG22	1:E:79:LEU:N	2.33	0.43
1:H:178:ARG:O	1:H:187:GLN:HB3	2.18	0.43
1:D:203:GLY:HA3	2:D:271:HOH:O	2.18	0.43
1:B:64:ARG:HD2	1:B:103:GLN:OE1	2.19	0.43
1:B:182:ASN:HB2	1:B:183:PRO:HD2	2.00	0.43
1:E:90:TYR:CE2	1:E:131:LYS:HD3	2.53	0.43
1:G:85:MET:HG3	1:G:245:MET:SD	2.58	0.43
1:C:65:GLY:O	1:C:74:LEU:HA	2.19	0.43
1:G:86:ASP:OD1	1:G:88:ARG:HB2	2.19	0.42
1:H:42:GLU:HG3	1:I:209:LEU:HD21	2.01	0.42
1:I:54:ALA:HB1	1:I:141:ILE:HD11	2.00	0.42
1:I:189:PHE:CE1	1:I:236:ILE:HD13	2.54	0.42
1:D:133:ASN:HB2	2:D:265:HOH:O	2.19	0.42
1:E:55:ALA:HB3	1:E:56:PRO:HD3	2.01	0.42
1:G:125:THR:O	1:G:126:ASP:HB3	2.19	0.42
1:H:196:ILE:HB	1:H:219:SER:HB3	2.01	0.42
1:C:177:GLY:O	1:C:178:ARG:HD3	2.19	0.42
1:G:78:THR:HG22	1:G:79:LEU:H	1.84	0.42
1:G:141:ILE:O	1:G:141:ILE:HG23	2.19	0.42
1:D:167:THR:HG23	1:E:174:SER:HB3	2.00	0.42
1:E:114:PHE:CE2	1:E:134:ALA:HB2	2.54	0.42
1:C:182:ASN:OD1	1:C:184:THR:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:PHE:CD1	1:D:135:THR:HG23	2.55	0.42
1:A:227:GLU:HB2	2:C:282:HOH:O	2.19	0.42
1:B:51:VAL:O	1:B:55:ALA:HB3	2.20	0.42
1:C:40:THR:HB	1:C:42:GLU:OE2	2.20	0.42
1:G:167:THR:HG23	1:H:174:SER:HB3	2.02	0.42
1:A:141:ILE:HG23	1:A:141:ILE:O	2.20	0.41
1:B:64:ARG:HD3	1:B:74:LEU:HB3	2.02	0.41
1:G:194:ALA:O	1:G:196:ILE:HG12	2.20	0.41
1:E:244:ILE:O	1:E:247:LYS:HG2	2.20	0.41
1:C:217:THR:HG22	1:C:218:LEU:H	1.84	0.41
1:F:244:ILE:O	1:F:248:LEU:HG	2.20	0.41
1:E:108:LEU:HD12	1:E:108:LEU:N	2.36	0.41
1:G:90:TYR:CE2	1:G:131:LYS:HD3	2.56	0.41
1:F:182:ASN:HB2	1:F:183:PRO:HD2	2.03	0.41
1:G:182:ASN:HB2	1:G:183:PRO:HD2	2.03	0.41
1:E:170:GLN:HB2	1:F:172:ILE:HD13	2.03	0.40
1:F:129:VAL:HG23	1:F:248:LEU:HD12	2.02	0.40
1:B:197:ASN:HB3	1:B:198:PRO:HD2	2.02	0.40
1:H:94:ASN:OD1	1:H:96:HIS:HB2	2.21	0.40
1:H:190:LEU:O	1:H:234:PHE:HA	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	198/241 (82%)	194 (98%)	4 (2%)	0	100	100
1	B	200/241 (83%)	197 (98%)	3 (2%)	0	100	100
1	C	199/241 (83%)	196 (98%)	3 (2%)	0	100	100
1	D	197/241 (82%)	193 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	187/241 (78%)	181 (97%)	6 (3%)	0	100	100
1	F	203/241 (84%)	198 (98%)	5 (2%)	0	100	100
1	G	194/241 (80%)	191 (98%)	3 (2%)	0	100	100
1	H	191/241 (79%)	186 (97%)	5 (3%)	0	100	100
1	I	185/241 (77%)	177 (96%)	8 (4%)	0	100	100
All	All	1754/2169 (81%)	1713 (98%)	41 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	162/198 (82%)	162 (100%)	0	100	100
1	B	165/198 (83%)	165 (100%)	0	100	100
1	C	164/198 (83%)	164 (100%)	0	100	100
1	D	162/198 (82%)	162 (100%)	0	100	100
1	E	155/198 (78%)	155 (100%)	0	100	100
1	F	164/198 (83%)	164 (100%)	0	100	100
1	G	158/198 (80%)	158 (100%)	0	100	100
1	H	153/198 (77%)	153 (100%)	0	100	100
1	I	152/198 (77%)	152 (100%)	0	100	100
All	All	1435/1782 (80%)	1435 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	188	ASN

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Mol	Chain	Res	Type
1	G	94	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	204/241 (84%)	0.42	14 (6%) 16 18	24, 45, 72, 85	0
1	B	206/241 (85%)	0.43	22 (10%) 6 5	26, 44, 71, 83	0
1	C	205/241 (85%)	0.41	16 (7%) 13 14	26, 46, 74, 83	0
1	D	203/241 (84%)	0.49	21 (10%) 6 6	30, 49, 71, 81	1 (0%)
1	E	195/241 (80%)	0.54	15 (7%) 13 15	34, 56, 75, 80	0
1	F	207/241 (85%)	0.36	11 (5%) 26 29	27, 43, 71, 82	0
1	G	200/241 (82%)	0.43	20 (10%) 7 7	38, 53, 75, 81	0
1	H	197/241 (81%)	0.58	22 (11%) 5 5	42, 60, 77, 82	0
1	I	193/241 (80%)	0.86	29 (15%) 2 2	43, 64, 80, 83	0
All	All	1810/2169 (83%)	0.50	170 (9%) 8 9	24, 52, 75, 85	1 (0%)

All (170) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	125	THR	7.4
1	D	39	SER	7.4
1	F	68	THR	6.5
1	F	69	ASN	6.5
1	H	40	THR	6.3
1	I	123	SER	5.7
1	A	225	ASP	5.6
1	D	40	THR	5.3
1	A	226	GLY	5.0
1	A	228	THR	5.0
1	E	221	ASP	4.9
1	B	67	ASN	4.9
1	E	188	ASN	4.8
1	H	77	ARG	4.7
1	H	220	PHE	4.7

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Mol	Chain	Res	Type	RSRZ
1	I	220	PHE	4.7
1	B	221	ASP	4.7
1	B	66	LEU	4.7
1	H	137	GLY	4.6
1	B	41	ASP	4.6
1	I	124	LEU	4.5
1	A	227	GLU	4.5
1	B	39	SER	4.5
1	I	219	SER	4.3
1	H	228	THR	4.3
1	I	244	ILE	4.3
1	I	74	LEU	4.2
1	F	221	ASP	4.2
1	G	247	LYS	4.2
1	H	248	LEU	4.2
1	B	219	SER	4.1
1	E	66	LEU	4.0
1	E	133	ASN	4.0
1	A	180	GLY	4.0
1	H	102	ASP	4.0
1	D	73	GLN	3.9
1	E	220	PHE	3.9
1	C	65	GLY	3.9
1	B	72	ASN	3.9
1	H	100	ASP	3.9
1	F	39	SER	3.8
1	F	43	THR	3.8
1	G	100	ASP	3.8
1	B	249	ILE	3.8
1	B	69	ASN	3.7
1	D	69	ASN	3.6
1	D	133	ASN	3.6
1	E	119	VAL	3.6
1	G	219	SER	3.6
1	A	221	ASP	3.5
1	I	146	ARG	3.5
1	C	74	LEU	3.5
1	H	63	ASN	3.5
1	G	228	THR	3.5
1	A	222	LYS	3.4
1	A	179	ILE	3.4
1	E	179	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	G	220	PHE	3.4
1	H	105	ILE	3.4
1	G	135	THR	3.4
1	B	74	LEU	3.3
1	G	250	ARG	3.3
1	I	136	GLY	3.3
1	G	40	THR	3.3
1	G	249	ILE	3.3
1	G	41	ASP	3.3
1	F	70	SER	3.2
1	B	220	PHE	3.2
1	D	70	SER	3.2
1	B	40	THR	3.2
1	C	75	GLU	3.2
1	A	220	PHE	3.1
1	F	41	ASP	3.1
1	I	236	ILE	3.1
1	B	38	ASP	3.1
1	A	66	LEU	3.1
1	D	65	GLY	3.1
1	B	68	THR	3.1
1	D	71	HIS	3.1
1	C	40	THR	3.1
1	A	188	ASN	3.1
1	H	227	GLU	3.0
1	G	39	SER	3.0
1	H	75	GLU	2.9
1	B	123	SER	2.9
1	D	219	SER	2.9
1	D	187	GLN	2.9
1	I	137	GLY	2.9
1	I	145	ALA	2.9
1	G	227	GLU	2.9
1	C	102	ASP	2.9
1	D	254	VAL	2.9
1	H	104	ILE	2.8
1	H	136	GLY	2.8
1	C	38	ASP	2.8
1	H	78	THR	2.8
1	E	171	GLY	2.8
1	I	138	LEU	2.8
1	G	43	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	H	135	THR	2.8
1	E	111	GLY	2.7
1	C	227	GLU	2.7
1	E	236	ILE	2.7
1	F	67	ASN	2.7
1	H	251	ASP	2.7
1	I	179	ILE	2.7
1	H	148	VAL	2.7
1	I	157	ALA	2.7
1	B	77	ARG	2.7
1	B	124	LEU	2.7
1	I	178	ARG	2.7
1	D	220	PHE	2.7
1	C	101	ALA	2.6
1	F	71	HIS	2.6
1	D	77	ARG	2.6
1	G	127	LEU	2.6
1	E	189	PHE	2.6
1	I	110	ASP	2.6
1	C	66	LEU	2.6
1	C	39	SER	2.5
1	C	248	LEU	2.5
1	G	77	ARG	2.5
1	F	42	GLU	2.5
1	I	101	ALA	2.5
1	D	66	LEU	2.5
1	B	73	GLN	2.5
1	C	250	ARG	2.5
1	D	41	ASP	2.5
1	I	240	LEU	2.5
1	I	77	ARG	2.5
1	G	38	ASP	2.5
1	C	77	ARG	2.4
1	C	228	THR	2.4
1	I	134	ALA	2.4
1	F	247	LYS	2.4
1	I	90	TYR	2.4
1	D	244	ILE	2.4
1	B	250	ARG	2.4
1	I	221	ASP	2.4
1	H	226	GLY	2.3
1	B	247	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	250	ARG	2.3
1	A	224	ASN	2.3
1	H	247	LYS	2.3
1	I	85	MET	2.3
1	H	106	VAL	2.2
1	I	229	PRO	2.2
1	G	171	GLY	2.2
1	E	222	LYS	2.2
1	C	119	VAL	2.1
1	D	218	LEU	2.1
1	E	134	ALA	2.1
1	I	112	ARG	2.1
1	E	114	PHE	2.1
1	G	119	VAL	2.1
1	G	63	ASN	2.1
1	D	102	ASP	2.1
1	H	146	ARG	2.1
1	E	243	LYS	2.1
1	G	170	GLN	2.1
1	A	135	THR	2.1
1	I	135	THR	2.1
1	B	42	GLU	2.1
1	I	96	HIS	2.1
1	I	169	THR	2.1
1	D	127	LEU	2.0
1	C	117	LEU	2.0
1	D	148	VAL	2.0
1	B	43	THR	2.0
1	A	223	SER	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.