



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2017 – 09:49 AM EDT

PDB ID : 3GDV
Title : Crystal structure of DegS H198P/D320A mutant modified by DFP and in complex with YQF peptide
Authors : Sohn, J.; Grant, R.A.; Sauer, R.T.
Deposited on : unknown
Resolution : 2.49 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

<http://wwpdb.org/validation/2016/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
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

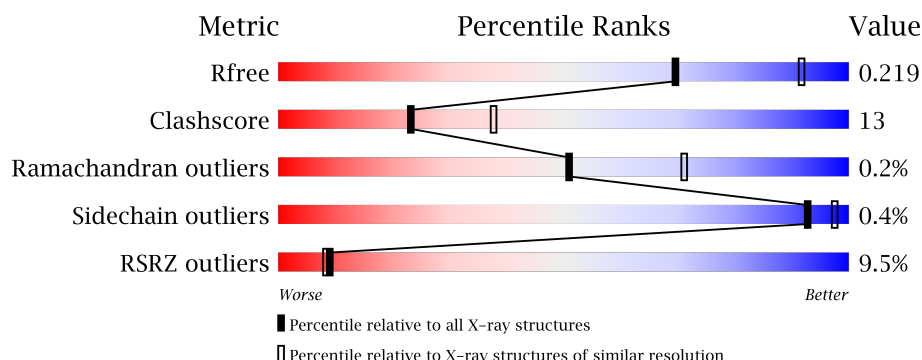
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	100719	4719 (2.50-2.46)
Clashscore	112137	5483 (2.50-2.46)
Ramachandran outliers	110173	5388 (2.50-2.46)
Sidechain outliers	110143	5390 (2.50-2.46)
RSRZ outliers	101464	4754 (2.50-2.46)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	340	<div> <div>6%</div> <div>65%</div> <div>22%</div> <div>13%</div> </div>
1	B	340	<div> <div>14%</div> <div>66%</div> <div>18%</div> <div>16%</div> </div>
1	C	340	<div> <div>4%</div> <div>69%</div> <div>17%</div> <div>14%</div> </div>
2	D	3	<div> <div>67%</div> <div>67%</div> <div>33%</div> </div>
2	E	3	<div> <div>100%</div> <div>67%</div> <div>33%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	3	 A horizontal bar chart showing the quality of chain F. The bar is divided into two segments: a green segment on the left representing 33% and a yellow segment on the right representing 67%. The percentages are labeled below the bar.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6662 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 DegS protease.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	P	S	0	0	0
			2193	1371	388	428	1	5			
1	B	284	Total	C	N	O	P	S	0	0	0
			2103	1321	367	410	1	4			
1	C	294	Total	C	N	O	P	S	0	0	0
			2189	1372	388	423	1	5			

There are 39 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
A	198	PRO	HIS	ENGINEERED	UNP P0AEE3
A	320	ALA	ASP	ENGINEERED	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

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Chain	Residue	Modelled	Actual	Comment	Reference
B	26	GLY	-	EXPRESSION TAG	UNP P0AEE3
B	198	PRO	HIS	ENGINEERED	UNP P0AEE3
B	320	ALA	ASP	ENGINEERED	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	UNP P0AEE3
C	320	ALA	ASP	ENGINEERED	UNP P0AEE3

- Molecule 2 is a protein called YQF peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			33	23	4	6			
2	E	3	Total	C	N	O	0	0	0
			33	23	4	6			
2	F	3	Total	C	N	O	0	0	0
			33	23	4	6			

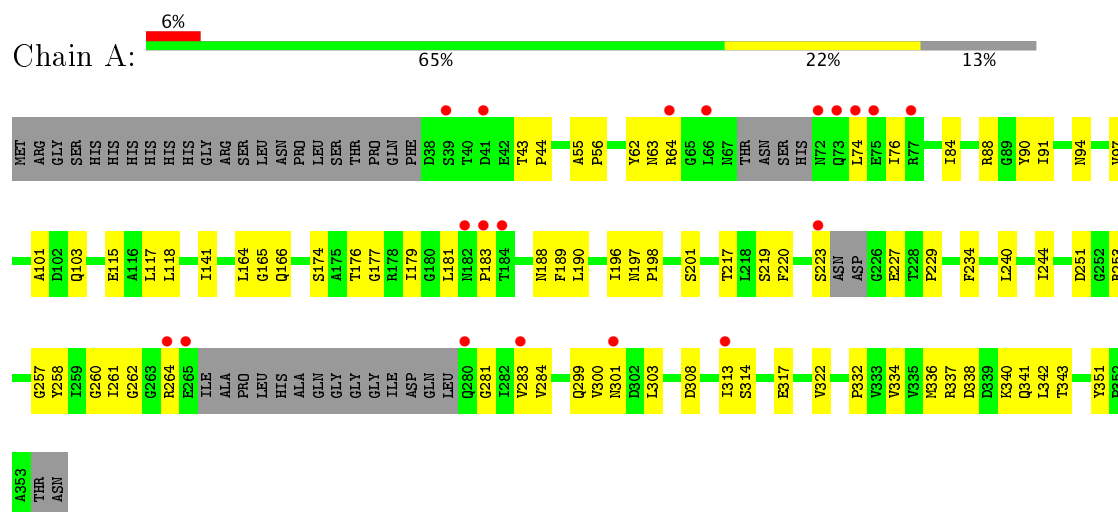
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		
3	B	34	Total	O	0	0
			34	34		
3	C	32	Total	O	0	0
			32	32		

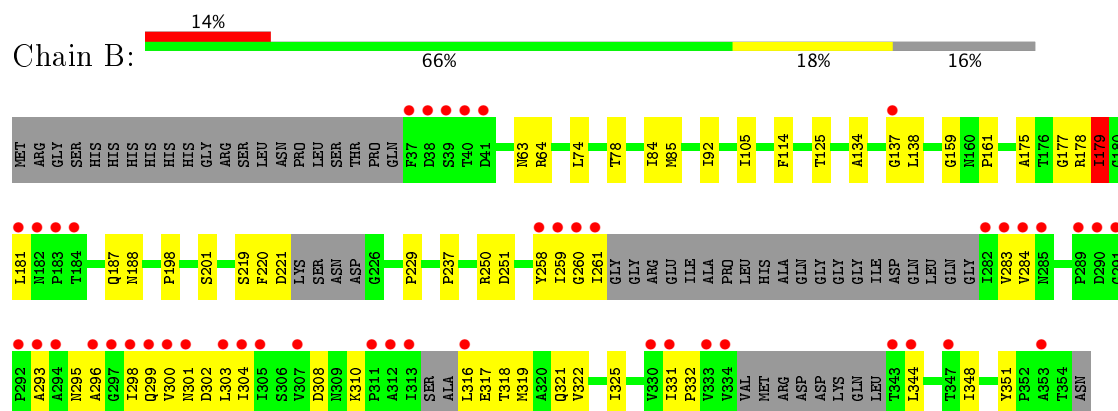
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

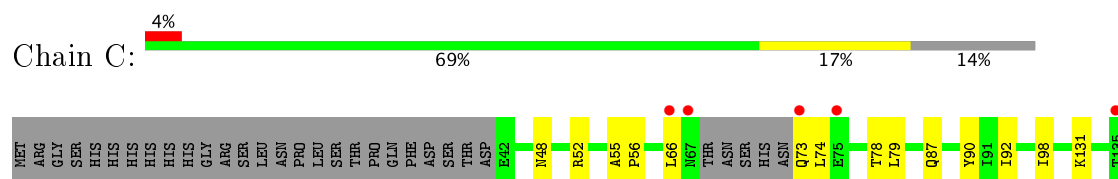
• Molecule 1: DegS protease

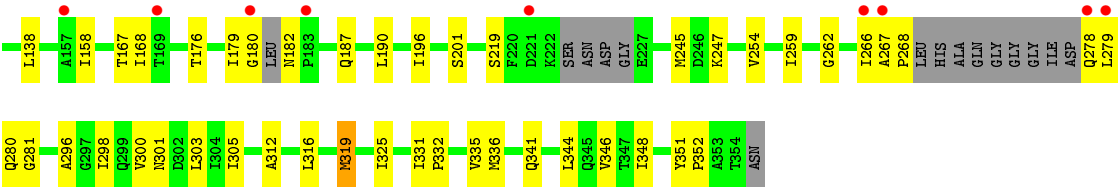


• Molecule 1: DegS protease



• Molecule 1: DegS protease





● Molecule 2: YQF peptide



● Molecule 2: YQF peptide



● Molecule 2: YQF peptide



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.88Å 172.28Å 114.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.04 – 2.49 32.04 – 2.49	Depositor EDS
% Data completeness (in resolution range)	95.9 (32.04-2.49) 95.9 (32.04-2.49)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 2.48Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.191 , 0.224 0.184 , 0.219	Depositor DCC
R_{free} test set	2001 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 63.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6662	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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: MIS

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	0.24	0/2204	0.45	0/2996
1	B	0.29	0/2115	0.50	0/2881
1	C	0.27	0/2200	0.49	0/2989
2	D	0.29	0/34	0.27	0/43
2	E	0.32	0/34	0.28	0/43
2	F	0.33	0/34	0.27	0/43
All	All	0.27	0/6621	0.48	0/8995

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	2193	0	2232	63	0
1	B	2103	0	2129	66	0
1	C	2189	0	2239	49	0
2	D	33	0	25	1	0
2	E	33	0	25	2	0
2	F	33	0	25	5	0
3	A	12	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	34	0	0	0	0
3	C	32	0	0	0	0
All	All	6662	0	6675	170	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 13.

All (170) 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:B:261:ILE:HG13	1:B:293:ALA:HB1	1.45	0.96
1:B:181:LEU:HD21	1:B:220:PHE:CD1	2.05	0.91
1:C:87:GLN:HE22	1:C:138:LEU:H	1.16	0.90
1:B:178:ARG:O	1:B:179:ILE:HG12	1.72	0.89
1:A:179:ILE:HD12	1:A:183:PRO:HA	1.57	0.87
1:B:258:TYR:HB2	1:B:351:TYR:HB2	1.54	0.86
1:A:62:TYR:HB3	1:A:64:ARG:HH12	1.43	0.82
1:B:261:ILE:HG13	1:B:293:ALA:CB	2.10	0.81
1:B:298:ILE:HG12	1:B:344:LEU:HD13	1.62	0.81
1:A:181:LEU:CD2	1:A:220:PHE:HD1	1.97	0.76
1:A:334:VAL:HG22	1:A:343:THR:HG22	1.67	0.76
1:A:64:ARG:CZ	1:A:76:ILE:HG12	2.17	0.75
1:B:84:ILE:HG21	1:B:138:LEU:HD13	1.68	0.75
1:A:74:LEU:HD23	1:A:74:LEU:H	1.52	0.75
1:C:196:ILE:CG2	1:C:201:MIS:H31	2.18	0.74
1:A:223:SER:HA	1:A:227:GLU:HB2	1.71	0.72
1:A:62:TYR:HB3	1:A:64:ARG:NH1	2.03	0.72
1:C:196:ILE:HG22	1:C:201:MIS:H21	1.71	0.71
1:B:181:LEU:CD2	1:B:220:PHE:CD1	2.74	0.71
1:B:201:MIS:H21	1:B:219:SER:HB2	1.73	0.70
1:C:73:GLN:HG2	1:C:74:LEU:H	1.59	0.67
1:B:261:ILE:CG1	1:B:293:ALA:HB1	2.24	0.66
1:A:117:LEU:HD12	1:A:118:LEU:H	1.61	0.65
1:A:181:LEU:HD22	1:A:220:PHE:HD1	1.62	0.64
1:B:296:ALA:O	1:B:344:LEU:HD22	1.97	0.64
1:C:87:GLN:NE2	1:C:138:LEU:H	1.91	0.64
1:B:179:ILE:HA	1:B:187:GLN:O	1.99	0.63
1:B:181:LEU:HD21	1:B:220:PHE:HD1	1.61	0.62
1:B:64:ARG:HH21	1:B:74:LEU:HD13	1.62	0.62
1:C:316:LEU:HD21	2:F:408:TYR:HE1	1.63	0.62
1:A:164:LEU:O	1:B:178:ARG:NH1	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:337:ARG:O	1:A:340:LYS:HG2	1.99	0.62
1:C:266:ILE:HD13	1:C:278:GLN:NE2	2.15	0.61
1:C:259:ILE:HD11	2:F:410:PHE:CE2	2.35	0.61
1:C:180:GLY:HA3	1:C:182:ASN:N	2.15	0.61
1:B:317:GLU:HA	1:B:317:GLU:OE1	2.00	0.61
1:B:181:LEU:HD11	1:B:220:PHE:HA	1.83	0.61
1:A:164:LEU:C	1:B:178:ARG:HH12	2.04	0.60
1:A:308:ASP:HA	1:A:332:PRO:HD2	1.83	0.60
1:B:308:ASP:HB2	1:B:331:ILE:HD13	1.83	0.60
1:B:322:VAL:HA	1:B:325:ILE:HD12	1.82	0.59
1:A:94:ASN:HD21	1:A:217:THR:HA	1.67	0.59
1:B:258:TYR:CZ	1:B:260:GLY:HA2	2.38	0.58
1:A:64:ARG:NH1	1:A:76:ILE:HG12	2.18	0.58
1:A:336:MET:HA	1:A:340:LYS:O	2.04	0.58
1:C:298:ILE:HG12	1:C:344:LEU:HD12	1.85	0.58
1:C:87:GLN:HE22	1:C:138:LEU:N	1.96	0.58
1:B:319:MET:HA	1:B:322:VAL:HG12	1.86	0.57
1:A:181:LEU:HD22	1:A:220:PHE:CD1	2.39	0.57
1:B:198:PRO:HA	1:B:201:MIS:H33	1.86	0.57
1:B:325:ILE:HD13	1:B:331:ILE:CD1	2.35	0.57
1:C:73:GLN:CG	1:C:74:LEU:H	2.16	0.57
1:B:299:GLN:O	1:B:302:ASP:HB2	2.05	0.56
1:B:300:VAL:O	1:B:301:ASN:HB2	2.06	0.56
1:C:268:PRO:HA	1:C:280:GLN:HE22	1.70	0.55
1:B:250:ARG:HG2	1:B:251:ASP:OD1	2.07	0.55
1:A:117:LEU:HD12	1:A:118:LEU:N	2.21	0.55
1:A:258:TYR:CZ	1:A:260:GLY:HA2	2.42	0.55
1:A:300:VAL:O	1:A:301:ASN:HB2	2.07	0.55
1:B:298:ILE:HG12	1:B:344:LEU:CD1	2.34	0.55
1:A:196:ILE:HG21	1:A:201:MIS:H31	1.88	0.54
1:A:84:ILE:HD13	1:A:91:ILE:HG12	1.88	0.54
1:B:261:ILE:HD11	1:B:298:ILE:HD12	1.90	0.54
1:A:165:GLY:CA	1:B:178:ARG:HH12	2.21	0.54
1:A:164:LEU:C	1:B:178:ARG:NH1	2.61	0.53
1:A:177:GLY:HA2	1:A:188:ASN:OD1	2.08	0.53
1:A:196:ILE:CG2	1:A:201:MIS:H31	2.37	0.53
1:C:92:ILE:HD11	1:C:245:MET:HB2	1.90	0.53
1:A:314:SER:HB3	1:A:317:GLU:OE1	2.08	0.53
1:A:64:ARG:HD2	1:A:74:LEU:HB2	1.90	0.53
1:B:259:ILE:HG22	2:E:410:PHE:O	2.07	0.53
1:C:196:ILE:HG21	1:C:201:MIS:H31	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:325:ILE:HG21	1:C:348:ILE:HG13	1.91	0.52
1:A:74:LEU:H	1:A:74:LEU:CD2	2.19	0.52
1:C:279:LEU:HD13	1:C:305:ILE:CD1	2.40	0.52
1:C:66:LEU:HA	1:C:73:GLN:O	2.09	0.52
1:B:319:MET:HA	1:B:322:VAL:CG1	2.40	0.52
1:B:295:ASN:O	1:B:296:ALA:HB3	2.10	0.52
1:C:90:TYR:CE2	1:C:131:LYS:HD3	2.45	0.51
1:A:174:SER:HB3	1:C:167:THR:HG23	1.91	0.51
1:B:125:THR:HA	1:B:187:GLN:NE2	2.25	0.51
1:B:64:ARG:NH2	1:B:74:LEU:HD13	2.25	0.51
1:A:181:LEU:HD21	1:A:220:PHE:HD1	1.73	0.51
1:C:247:LYS:HD2	1:C:254:VAL:HG13	1.93	0.51
1:A:299:GLN:CD	1:A:337:ARG:HH12	2.13	0.51
1:C:279:LEU:HD13	1:C:305:ILE:HD11	1.93	0.51
1:B:316:LEU:O	1:B:319:MET:HG2	2.10	0.51
1:A:88:ARG:HD2	1:A:90:TYR:HE1	1.75	0.50
1:B:198:PRO:CA	1:B:201:MIS:H33	2.42	0.50
1:C:196:ILE:HG23	1:C:201:MIS:H31	1.93	0.50
1:A:261:ILE:HG23	1:A:284:VAL:HG13	1.94	0.50
1:A:220:PHE:HD2	1:A:229:PRO:CB	2.25	0.50
1:B:299:GLN:HG2	1:B:300:VAL:O	2.11	0.50
1:B:283:VAL:HG22	1:B:303:LEU:HD23	1.94	0.49
1:B:318:THR:O	1:B:322:VAL:HG12	2.11	0.49
1:A:251:ASP:C	1:A:253:ARG:H	2.16	0.49
1:C:296:ALA:HB3	1:C:346:VAL:HG12	1.94	0.49
1:A:264:ARG:O	1:A:283:VAL:HG12	2.12	0.49
1:B:310:LYS:HG3	1:B:321:GLN:NE2	2.27	0.49
1:C:201:MIS:H22	1:C:219:SER:HB2	1.94	0.49
1:A:281:GLY:HA3	1:A:303:LEU:HD11	1.93	0.49
1:B:259:ILE:HG22	2:E:410:PHE:C	2.33	0.49
1:C:78:THR:HG22	1:C:79:LEU:N	2.28	0.49
1:A:341:GLN:O	1:A:342:LEU:HD23	2.13	0.48
1:A:166:GLN:O	1:B:178:ARG:NH2	2.47	0.48
1:C:319:MET:HE2	2:F:410:PHE:HB3	1.96	0.48
1:A:220:PHE:HD2	1:A:229:PRO:HB3	1.79	0.48
1:A:165:GLY:C	1:B:178:ARG:HH22	2.17	0.48
1:A:262:GLY:HA2	2:D:408:TYR:O	2.14	0.47
1:B:303:LEU:HD13	1:B:304:ILE:N	2.29	0.47
1:A:299:GLN:CG	1:A:337:ARG:HH12	2.27	0.47
1:B:177:GLY:HA2	1:B:188:ASN:OD1	2.14	0.47
1:B:178:ARG:O	1:B:179:ILE:CG1	2.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:MET:HG3	1:B:92:ILE:HG13	1.97	0.47
1:A:258:TYR:HB2	1:A:351:TYR:HA	1.95	0.47
1:B:261:ILE:HD11	1:B:298:ILE:CD1	2.45	0.47
1:C:262:GLY:HA2	2:F:408:TYR:O	2.15	0.47
1:C:73:GLN:HG2	1:C:74:LEU:N	2.26	0.46
1:C:305:ILE:HD11	1:C:336:MET:HG2	1.96	0.46
1:A:257:GLY:HA3	1:A:322:VAL:O	2.16	0.46
1:A:74:LEU:N	1:A:74:LEU:HD23	2.24	0.46
1:B:114:PHE:CZ	1:B:134:ALA:HB2	2.51	0.46
1:B:331:ILE:HA	1:B:332:PRO:HD3	1.76	0.45
1:C:267:ALA:HB1	1:C:268:PRO:HD2	1.98	0.45
1:C:300:VAL:O	1:C:301:ASN:HB2	2.16	0.45
1:B:261:ILE:HG23	1:B:284:VAL:HG13	1.98	0.45
1:C:303:LEU:HB3	1:C:336:MET:HB2	1.99	0.45
1:A:338:ASP:HB2	1:A:340:LYS:HE3	1.99	0.45
1:C:158:ILE:HG23	1:C:168:ILE:HD13	1.98	0.45
1:B:188:ASN:O	1:B:237:PRO:HD3	2.17	0.44
1:B:64:ARG:HD3	1:B:105:ILE:HG12	2.00	0.44
1:B:325:ILE:HG21	1:B:348:ILE:HG13	1.99	0.44
1:A:176:THR:HG22	1:A:190:LEU:CD2	2.48	0.43
1:B:159:GLY:C	1:B:161:PRO:HD3	2.38	0.43
1:C:303:LEU:O	1:C:335:VAL:HA	2.18	0.43
1:A:240:LEU:O	1:A:244:ILE:HG12	2.18	0.43
1:B:293:ALA:O	1:B:298:ILE:HB	2.18	0.43
1:C:48:ASN:OD1	1:C:52:ARG:NH1	2.47	0.43
1:A:165:GLY:N	1:B:178:ARG:HH12	2.15	0.43
1:A:223:SER:CA	1:A:227:GLU:HB2	2.45	0.43
1:C:98:ILE:HD12	1:C:98:ILE:C	2.40	0.43
1:B:259:ILE:O	1:B:293:ALA:HB2	2.20	0.42
1:C:336:MET:SD	1:C:341:GLN:HG2	2.59	0.42
1:B:175:ALA:HB1	1:B:178:ARG:HE	1.84	0.42
1:B:259:ILE:CG2	1:B:261:ILE:H	2.32	0.42
1:C:55:ALA:N	1:C:56:PRO:CD	2.82	0.42
1:B:325:ILE:HD13	1:B:331:ILE:HD12	2.02	0.42
1:C:296:ALA:CB	1:C:346:VAL:HG12	2.48	0.42
1:C:316:LEU:CD2	2:F:408:TYR:HE1	2.31	0.42
1:C:305:ILE:HD11	1:C:336:MET:CG	2.51	0.41
1:A:63:ASN:ND2	1:A:101:ALA:HB2	2.35	0.41
1:A:84:ILE:CD1	1:A:91:ILE:HG12	2.49	0.41
1:B:259:ILE:O	1:B:293:ALA:N	2.54	0.41
1:A:141:ILE:O	1:A:141:ILE:HG23	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:351:TYR:HA	1:C:352:PRO:HD3	1.95	0.41
1:A:313:ILE:N	1:A:313:ILE:HD12	2.36	0.41
1:A:55:ALA:N	1:A:56:PRO:CD	2.83	0.41
1:A:94:ASN:HB2	1:A:97:VAL:HG23	2.02	0.41
1:A:43:THR:HA	1:A:44:PRO:HD3	1.89	0.41
1:C:176:THR:HG22	1:C:190:LEU:CD2	2.51	0.41
1:C:298:ILE:CG1	1:C:344:LEU:HD12	2.48	0.41
1:C:281:GLY:HA2	1:C:312:ALA:O	2.20	0.40
1:C:331:ILE:HA	1:C:332:PRO:HD3	1.95	0.40
1:A:103:GLN:HE21	1:A:115:GLU:CD	2.25	0.40
1:A:189:PHE:CG	1:A:234:PHE:HB3	2.57	0.40
1:A:197:ASN:HB3	1:A:198:PRO:HD2	2.04	0.40
1:B:221:ASP:HA	1:B:229:PRO:HG2	2.03	0.40
1:B:63:ASN:HB3	1:B:78:THR:HB	2.02	0.40
1:C:179:ILE:HD12	1:C:179:ILE:C	2.42	0.40
1:C:182:ASN:ND2	1:C:187:GLN:OE1	2.50	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	287/340 (84%)	279 (97%)	8 (3%)	0	100	100
1	B	273/340 (80%)	262 (96%)	9 (3%)	2 (1%)	25	41
1	C	283/340 (83%)	282 (100%)	1 (0%)	0	100	100
2	D	1/3 (33%)	1 (100%)	0	0	100	100
2	E	1/3 (33%)	1 (100%)	0	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100
All	All	846/1029 (82%)	826 (98%)	18 (2%)	2 (0%)	51	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	179	ILE
1	B	137	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	235/275 (86%)	234 (100%)	1 (0%)	93	97
1	B	226/275 (82%)	225 (100%)	1 (0%)	93	97
1	C	234/275 (85%)	233 (100%)	1 (0%)	93	97
2	D	3/3 (100%)	3 (100%)	0	100	100
2	E	3/3 (100%)	3 (100%)	0	100	100
2	F	3/3 (100%)	3 (100%)	0	100	100
All	All	704/834 (84%)	701 (100%)	3 (0%)	93	97

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

Mol	Chain	Res	Type
1	A	219	SER
1	B	179	ILE
1	C	319	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	94	ASN
1	A	103	GLN
1	A	295	ASN
1	A	321	GLN
1	B	187	GLN
1	B	295	ASN

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Mol	Chain	Res	Type
1	C	87	GLN
1	C	278	GLN
1	C	295	ASN
1	C	321	GLN
2	F	409	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	MIS	A	201	1	12,12,13	0.62	0	10,16,18	0.82	0
1	MIS	B	201	1	12,12,13	0.67	0	10,16,18	0.87	0
1	MIS	C	201	1	12,12,13	0.65	0	10,16,18	0.88	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	MIS	A	201	1	-	0/10/13/15	0/0/0/0
1	MIS	B	201	1	-	0/10/13/15	0/0/0/0
1	MIS	C	201	1	-	0/10/13/15	0/0/0/0

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 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	201	MIS	2	0
1	B	201	MIS	3	0
1	C	201	MIS	5	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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 ⓘ

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	295/340 (86%)	0.24	19 (6%)	20 20	39, 82, 143, 168	0
1	B	283/340 (83%)	0.57	46 (16%)	2 2	28, 50, 176, 198	0
1	C	293/340 (86%)	-0.03	14 (4%)	31 32	34, 54, 107, 150	0
2	D	3/3 (100%)	2.59	2 (66%)	0 0	128, 128, 139, 140	0
2	E	3/3 (100%)	3.55	3 (100%)	0 0	145, 145, 162, 164	0
2	F	3/3 (100%)	0.49	0	100 100	66, 66, 81, 99	0
All	All	880/1029 (85%)	0.27	84 (9%)	9 8	28, 60, 152, 198	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	ILE	11.1
1	B	313	ILE	8.3
1	A	66	LEU	6.9
1	B	181	LEU	6.5
1	A	74	LEU	6.4
1	B	291	GLY	6.0
1	B	282	ILE	5.9
1	B	301	ASN	5.9
1	B	289	PRO	5.8
1	B	316	LEU	5.6
1	B	182	ASN	5.4
1	A	264	ARG	5.3
1	B	37	PHE	5.2
2	E	408	TYR	5.2
1	B	305	ILE	5.1
1	B	297	GLY	5.0
1	B	303	LEU	4.9
1	B	183	PRO	4.6
1	B	294	ALA	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	261	ILE	4.5
1	B	300	VAL	4.5
1	B	283	VAL	4.1
1	A	313	ILE	4.0
2	D	410	PHE	4.0
1	A	73	GLN	4.0
1	B	258	TYR	3.9
1	A	301	ASN	3.8
1	B	299	GLN	3.8
1	B	344	LEU	3.8
1	B	296	ALA	3.7
1	A	280	GLN	3.6
1	A	72	ASN	3.5
1	B	293	ALA	3.5
1	B	38	ASP	3.5
1	C	267	ALA	3.5
1	C	73	GLN	3.4
1	B	312	ALA	3.3
1	C	266	ILE	3.3
1	B	284	VAL	3.3
1	B	285	ASN	3.2
1	C	75	GLU	3.2
1	C	183	PRO	3.2
2	D	408	TYR	3.1
1	B	39	SER	3.1
1	A	75	GLU	3.0
2	E	409	GLN	3.0
1	A	283	VAL	3.0
1	B	311	PRO	3.0
1	B	184	THR	2.9
1	A	184	THR	2.9
1	B	290	ASP	2.9
1	B	331	ILE	2.9
1	B	298	ILE	2.9
1	A	64	ARG	2.8
1	B	347	THR	2.8
1	C	135	THR	2.7
1	B	307	VAL	2.7
1	A	182	ASN	2.6
1	B	334	VAL	2.5
2	E	410	PHE	2.5
1	B	259	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	40	THR	2.4
1	B	41	ASP	2.4
1	B	292	PRO	2.3
1	B	260	GLY	2.3
1	A	265	GLU	2.3
1	C	157	ALA	2.3
1	B	330	VAL	2.2
1	A	183	PRO	2.2
1	A	77	ARG	2.2
1	B	353	ALA	2.2
1	A	41	ASP	2.2
1	C	180	GLY	2.2
1	C	279	LEU	2.2
1	B	343	THR	2.2
1	A	39	SER	2.1
1	B	333	VAL	2.1
1	C	66	LEU	2.1
1	C	169	THR	2.1
1	C	278	GLN	2.1
1	C	221	ASP	2.0
1	A	223	SER	2.0
1	C	67	ASN	2.0
1	B	137	GLY	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	MIS	C	201	13/14	0.99	0.18	-	30,34,46,54	0
1	MIS	A	201	13/14	0.99	0.12	-	40,51,60,67	0
1	MIS	B	201	13/14	0.98	0.15	-	26,35,60,78	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.