



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:11 AM UTC

PDB ID : 3GDV / pdb_00003gdv
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 : 2009-02-24
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

<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

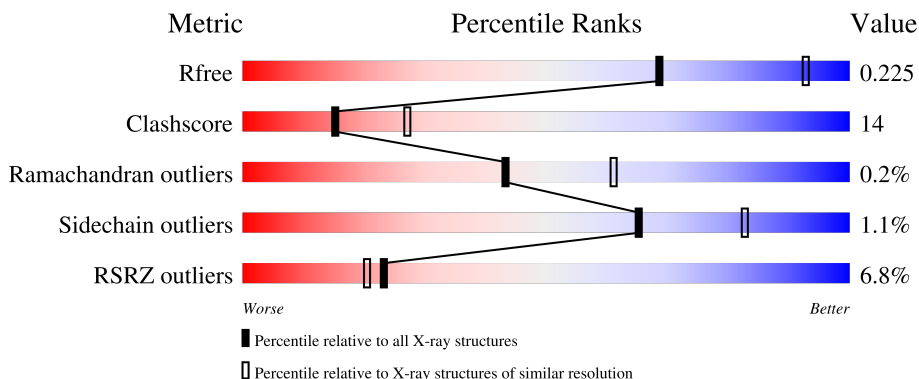
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	180053	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (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 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	340	3% 63% 24% 13%
1	B	340	10% 64% 19% 16%
1	C	340	3% 68% 18% 14%
2	D	3	33% 67% 33%
2	E	3	67% 33%

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Mol	Chain	Length	Quality of chain
2	F	3	33%33%67%

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 mutation	UNP P0AEE3
A	320	ALA	ASP	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

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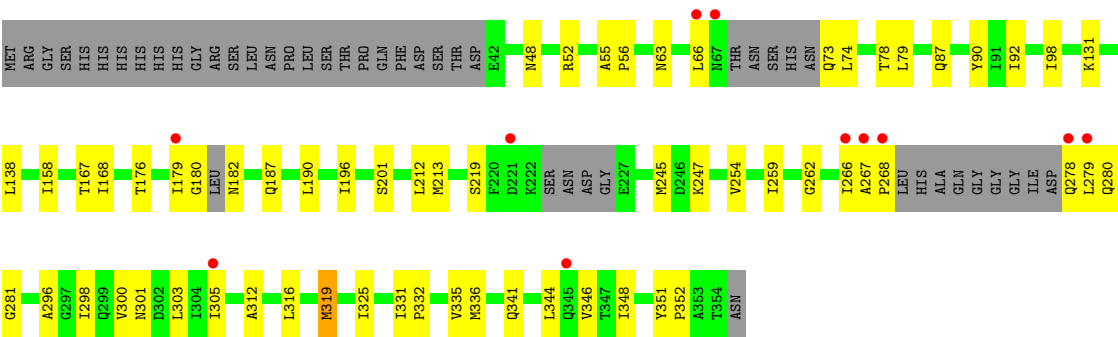
Chain	Residue	Modelled	Actual	Comment	Reference
B	26	GLY	-	expression tag	UNP P0AEE3
B	198	PRO	HIS	engineered mutation	UNP P0AEE3
B	320	ALA	ASP	engineered mutation	UNP P0AEE3
C	16	MET	-	expression tag	UNP P0AEE3
C	17	ARG	-	expression tag	UNP P0AEE3
C	18	GLY	-	expression tag	UNP P0AEE3
C	19	SER	-	expression 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
C	320	ALA	ASP	engineered mutation	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		



● 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) 95.8 (32.04-2.49)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.191 , 0.224 0.191 , 0.225	Depositor DCC
R_{free} test set	2003 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 72.0	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.32	0/2204	0.75	0/2996
1	B	0.39	0/2115	0.79	1/2881 (0.0%)
1	C	0.36	0/2200	0.76	0/2989
2	D	0.26	0/34	0.35	0/43
2	E	0.18	0/34	0.34	0/43
2	F	0.17	0/34	0.37	0/43
All	All	0.36	0/6621	0.76	1/8995 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	184	THR	CB-CA-C	-6.03	109.61	116.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2193	0	2232	66	0
1	B	2103	0	2129	69	0
1	C	2189	0	2239	53	0
2	D	33	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	33	0	25	3	0
2	F	33	0	25	5	0
3	A	12	0	0	0	0
3	B	34	0	0	0	0
3	C	32	0	0	0	0
All	All	6662	0	6675	180	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 14.

All (180) 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:H	1:A:74:LEU:HD23	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:164:LEU:C	1:B:178:ARG:HH12	2.04	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:ARG:HH21	1:B:74:LEU:HD13	1.62	0.62
1:B:181:LEU:HD21	1:B:220:PHE:HD1	1.61	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
1:A:337:ARG:O	1:A:340:LYS:HG2	1.99	0.62
1:C:259:ILE:HD11	2:F:410:PHE:CE2	2.35	0.61
1:C:266:ILE:HD13	1:C:278:GLN:NE2	2.15	0.61
1:C:180:GLY:HA3	1:C:182:ASN:N	2.15	0.61
1:B:181:LEU:HD11	1:B:220:PHE:HA	1.83	0.61
1:B:317:GLU:HA	1:B:317:GLU:OE1	2.00	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:A:207:ASN:HB3	1:A:213:MET:HE2	1.84	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:A:164:LEU:C	1:B:178:ARG:NH1	2.61	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:A:299:GLN:CD	1:A:337:ARG:HH12	2.13	0.55
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:B:213:MET:HA	1:B:213:MET:HE2	1.87	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:B:259:ILE:HG22	2:E:410:PHE:C	2.33	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:251:ASP:C	1:A:253:ARG:H	2.16	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
1:A:165:GLY:C	1:B:178:ARG:HH22	2.17	0.52
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:66:LEU:HA	1:C:73:GLN:O	2.09	0.52
1:C:279:LEU:HD13	1:C:305:ILE:CD1	2.40	0.52
1:B:85:MET:HE2	1:B:92:ILE:HD12	1.92	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:A:174:SER:HB3	1:C:167:THR:HG23	1.91	0.51
1:C:90:TYR:CE2	1:C:131:LYS:HD3	2.45	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: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:C:319:MET:HE2	2:F:410:PHE:HB3	1.94	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: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: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:B:159:GLY:C	1:B:161:PRO:HD3	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:GLN:O	1:B:178:ARG:NH2	2.47	0.48
1:C:212:LEU:O	1:C:213:MET:HE2	2.13	0.48
1:A:220:PHE:HD2	1:A:229:PRO:HB3	1.79	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:C:98:ILE:C	1:C:98:ILE:HD12	2.40	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
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:74:LEU:HD23	1:A:74:LEU:N	2.24	0.46
1:A:257:GLY:HA3	1:A:322:VAL:O	2.16	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:B:261:ILE:HG23	1:B:284:VAL:HG13	1.98	0.45
1:C:213:MET:HE2	1:C:213:MET:HA	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:179:ILE:C	1:C:179:ILE:HD12	2.42	0.45
1:A:103:GLN:HE21	1:A:115:GLU:CD	2.25	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:C:300:VAL:O	1:C:301:ASN:HB2	2.16	0.44
1:A:176:THR:HG22	1:A:190:LEU:CD2	2.48	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:63:ASN:OD1	1:C:63:ASN:C	2.63	0.42
1:A:79:LEU:C	1:A:79:LEU:HD12	2.45	0.41
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:C:179:ILE:O	1:C:180:GLY:C	2.63	0.41
1:A:141:ILE:O	1:A:141:ILE:HG23	2.19	0.41
1:C:351:TYR:HA	1:C:352:PRO:HD3	1.95	0.41
1:A:43:THR:HA	1:A:44:PRO:HD3	1.89	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:313:ILE:HD12	1:A:313:ILE:N	2.36	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: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:286:GLU:CD	2:E:409:GLN:HE21	2.29	0.40
1:B:63:ASN:HB3	1:B:78:THR:HB	2.02	0.40
1:B:221:ASP:HA	1:B:229:PRO:HG2	2.03	0.40
1:A:159:GLY:C	1:A:161:PRO:HD3	2.47	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 ⓘ

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	287/340 (84%)	279 (97%)	8 (3%)	0	100	100
1	B	273/340 (80%)	262 (96%)	9 (3%)	2 (1%)	18	32
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%)	43	61

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%)	231 (98%)	4 (2%)	53	75
1	B	226/275 (82%)	223 (99%)	3 (1%)	61	80
1	C	234/275 (85%)	233 (100%)	1 (0%)	84	92
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%)	696 (99%)	8 (1%)	65	83

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	SER
1	A	300	VAL
1	A	318	THR

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Mol	Chain	Res	Type
1	A	335	VAL
1	B	179	ILE
1	B	196	ILE
1	B	259	ILE
1	C	319	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	94	ASN
1	A	99	ASN
1	A	103	GLN
1	A	295	ASN
1	A	321	GLN
1	B	71	HIS
1	B	187	GLN
1	B	295	ASN
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	C	201	1	11,12,13	0.74	0	11,16,18	0.83	0
1	MIS	B	201	1	11,12,13	0.77	0	11,16,18	0.85	0
1	MIS	A	201	1	11,12,13	0.71	0	11,16,18	0.67	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	C	201	1	-	3/11/13/15	-
1	MIS	B	201	1	-	8/11/13/15	-
1	MIS	A	201	1	-	3/11/13/15	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	201	MIS	CB-OG-P-O2P
1	A	201	MIS	CB-OG-P-O3P
1	B	201	MIS	CB-OG-P-O1P
1	B	201	MIS	CB-OG-P-O2P
1	B	201	MIS	CB-OG-P-O3P
1	B	201	MIS	C1-O3P-P-OG
1	B	201	MIS	C3-C1-O3P-P
1	C	201	MIS	CB-OG-P-O3P
1	A	201	MIS	CA-CB-OG-P
1	B	201	MIS	CA-CB-OG-P
1	C	201	MIS	CA-CB-OG-P
1	C	201	MIS	CB-OG-P-O2P
1	B	201	MIS	C1-O3P-P-O1P
1	B	201	MIS	C1-O3P-P-O2P

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	201	MIS	5	0

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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%)	0.22	10 (3%) 48 44	39, 82, 143, 168	0
1	B	283/340 (83%)	0.14	35 (12%) 8 6	28, 50, 176, 198	0
1	C	293/340 (86%)	-0.15	11 (3%) 44 40	34, 54, 107, 150	0
2	D	3/3 (100%)	1.67	1 (33%) 1 1	128, 128, 139, 140	0
2	E	3/3 (100%)	1.88	2 (66%) 0 0	145, 145, 162, 164	0
2	F	3/3 (100%)	0.89	1 (33%) 1 1	66, 66, 81, 99	0
All	All	880/1029 (85%)	0.08	60 (6%) 23 20	28, 60, 152, 198	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	ILE	4.7
1	B	258	TYR	4.4
1	A	353	ALA	4.1
1	B	137	GLY	4.0
1	B	282	ILE	3.8
1	B	353	ALA	3.7
1	A	74	LEU	3.7
1	C	279	LEU	3.4
1	B	289	PRO	3.4
1	B	187	GLN	3.3
1	B	181	LEU	3.3
1	B	334	VAL	3.2
1	B	305	ILE	3.2
1	B	303	LEU	3.2
1	B	221	ASP	3.1
1	B	300	VAL	3.1
1	A	185	GLY	3.0
1	B	284	VAL	3.0
1	B	313	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	184	THR	3.0
1	C	278	GLN	3.0
1	C	67	ASN	3.0
1	A	223	SER	2.9
1	C	345	GLN	2.9
1	B	261	ILE	2.8
1	A	183	PRO	2.8
1	C	221	ASP	2.8
1	B	179	ILE	2.8
1	B	291	GLY	2.7
1	B	311	PRO	2.7
1	B	136	GLY	2.7
1	C	305	ILE	2.7
1	B	298	ILE	2.7
1	A	222	LYS	2.6
1	A	342	LEU	2.6
2	E	410	PHE	2.6
1	C	268	PRO	2.5
1	B	39	SER	2.5
1	A	64	ARG	2.5
1	C	266	ILE	2.5
1	B	135	THR	2.5
1	B	285	ASN	2.5
1	B	226	GLY	2.4
1	C	66	LEU	2.4
1	A	184	THR	2.4
1	B	345	GLN	2.4
2	F	408	TYR	2.3
2	E	408	TYR	2.3
1	A	186	ARG	2.3
1	B	352	PRO	2.3
1	B	301	ASN	2.2
1	B	283	VAL	2.2
1	B	185	GLY	2.2
1	C	267	ALA	2.2
1	B	287	VAL	2.1
1	B	294	ALA	2.1
1	B	292	PRO	2.1
1	B	227	GLU	2.1
2	D	408	TYR	2.1
1	C	179	ILE	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(Å ²)	Q<0.9
1	MIS	A	201	13/14	0.98	0.06	40,51,60,67	0
1	MIS	B	201	13/14	0.98	0.07	26,35,60,78	0
1	MIS	C	201	13/14	0.98	0.06	30,34,46,54	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.