



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:08 AM UTC

PDB ID : 3GDU / pdb_00003gdu
Title : Crystal structure of DegS H198P/D320A mutant modified by DFP and in complex with YRF peptide
Authors : Sohn, J.; Grant, R.A.; Sauer, R.T.
Deposited on : 2009-02-24
Resolution : 3.00 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

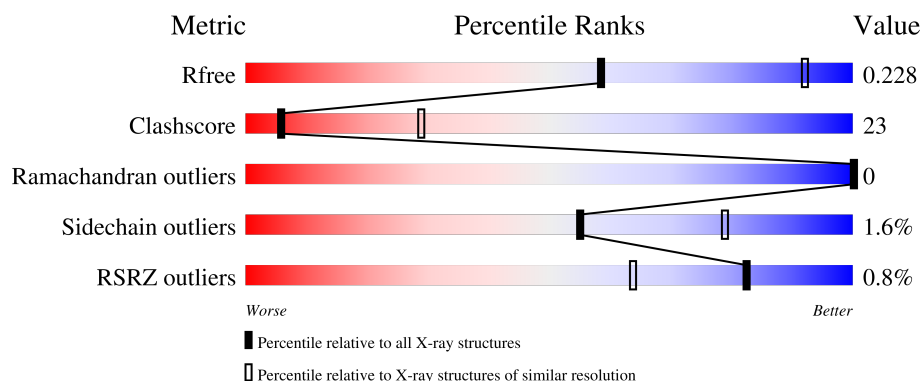
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	46% 39% 13%
1	B	340	54% 28% 18%
1	C	340	59% 27% 14%
2	D	3	67% 33%
2	E	3	67% 33%

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Mol	Chain	Length	Quality of chain
2	F	3	 100%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6530 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	295	Total	C	N	O	P	S	0	0	0
			2171	1355	378	432	1	5			
1	B	279	Total	C	N	O	P	S	0	0	0
			2069	1293	361	410	1	4			
1	C	293	Total	C	N	O	P	S	0	0	0
			2185	1366	388	425	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

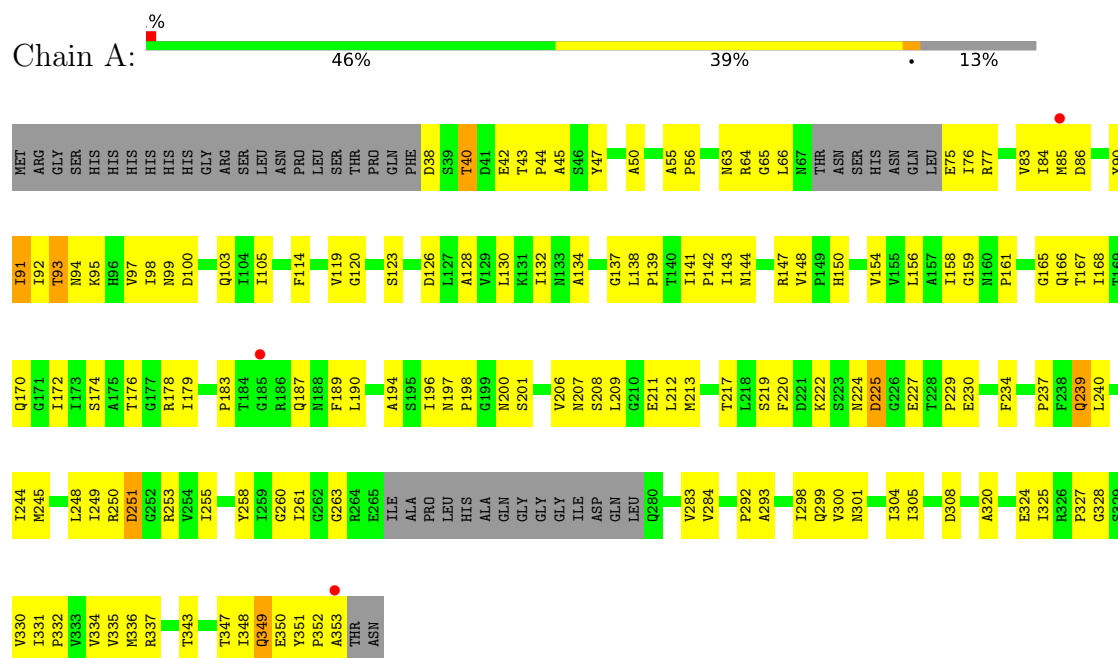
- Molecule 2 is a protein called YRF peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			35	24	6	5			
2	E	3	Total	C	N	O	0	0	0
			35	24	6	5			
2	F	3	Total	C	N	O	0	0	0
			35	24	6	5			

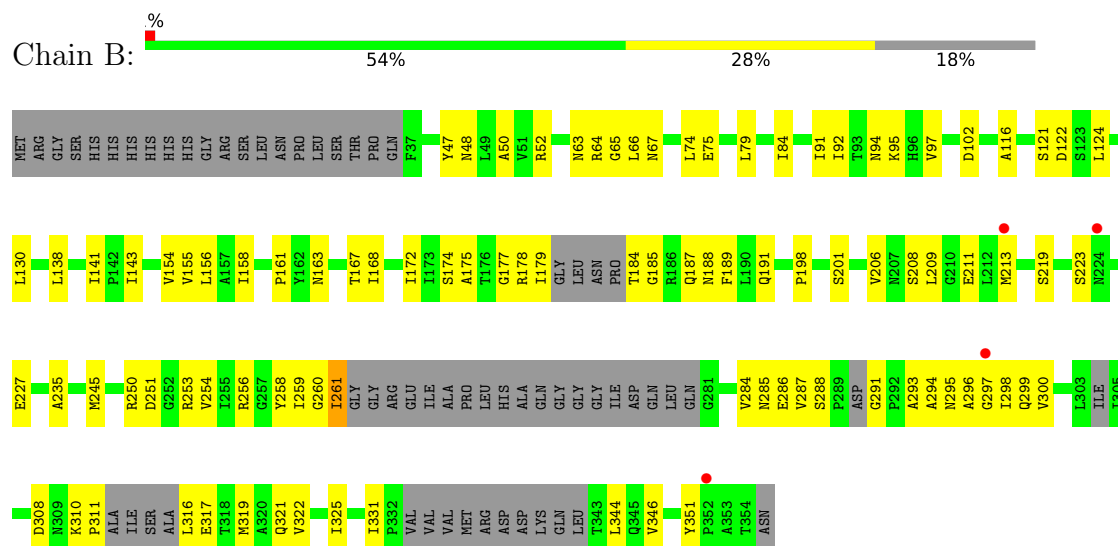
3 Residue-property plots

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.

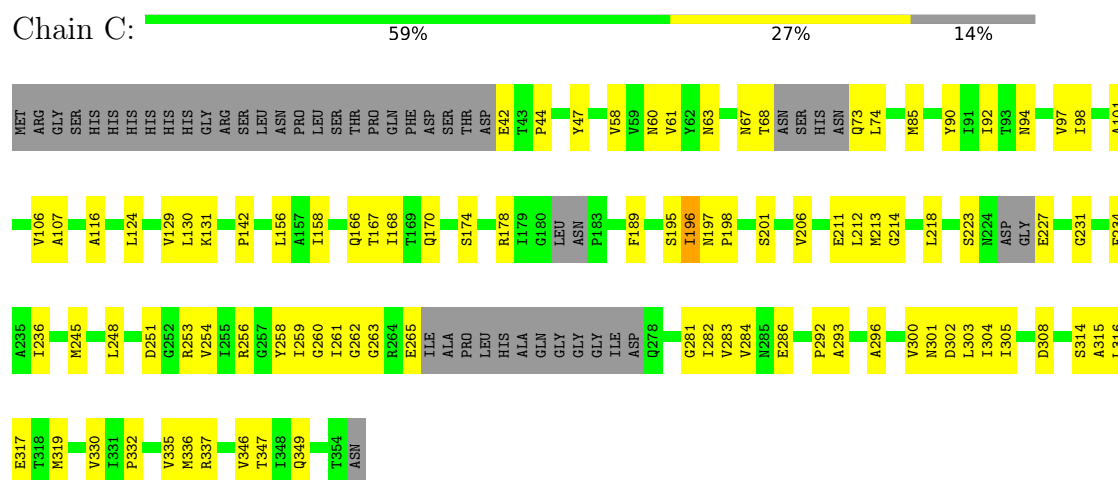
• Molecule 1: DegS protease



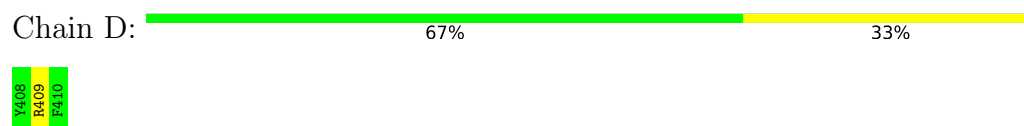
• Molecule 1: DegS protease



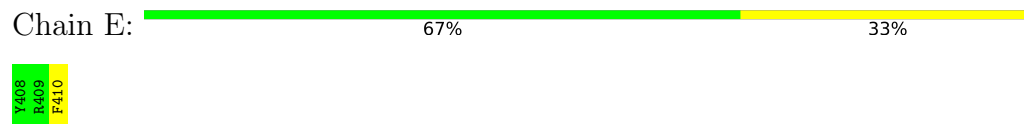
- Molecule 1: DegS protease



- Molecule 2: YRF peptide



- Molecule 2: YRF peptide



- Molecule 2: YRF peptide



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	117.57Å 171.28Å 111.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.15 – 3.00 36.15 – 3.00	Depositor EDS
% Data completeness (in resolution range)	92.9 (36.15-3.00) 94.9 (36.15-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.209 , 0.231 0.202 , 0.228	Depositor DCC
R_{free} test set	1117 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 55.6	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6530	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage'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 [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	0.28	0/2183	0.75	0/2974
1	B	0.29	0/2079	0.71	0/2828
1	C	0.27	0/2195	0.73	0/2980
2	D	0.30	0/36	0.41	0/45
2	E	0.20	0/36	0.46	0/45
2	F	0.13	0/36	0.37	0/45
All	All	0.28	0/6565	0.73	0/8917

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	2171	0	2180	140	0
1	B	2069	0	2075	96	0
1	C	2185	0	2233	81	0
2	D	35	0	30	1	0
2	E	35	0	30	2	0
2	F	35	0	30	3	0
All	All	6530	0	6578	301	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 23.

All (301) 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:A:305:ILE:HD11	1:A:336:MET:HG2	1.40	1.04
1:A:292:PRO:HG3	1:A:349:GLN:HG2	1.42	0.96
1:B:64:ARG:HH21	1:B:74:LEU:HD13	1.31	0.94
1:A:179:ILE:HD12	1:A:183:PRO:HA	1.57	0.85
1:A:224:ASN:O	1:A:227:GLU:HG2	1.78	0.84
1:B:258:TYR:HB2	1:B:351:TYR:HB2	1.63	0.81
1:A:174:SER:HB3	1:C:167:THR:HG23	1.64	0.80
1:B:293:ALA:O	1:B:298:ILE:HG22	1.82	0.80
1:B:178:ARG:O	1:B:179:ILE:HG12	1.83	0.79
1:A:299:GLN:HB2	1:A:337:ARG:HH12	1.47	0.78
1:A:85:MET:HE3	1:A:143:ILE:HD11	1.64	0.78
1:C:259:ILE:HD12	1:C:261:ILE:HD12	1.66	0.78
1:A:334:VAL:HG22	1:A:343:THR:HG22	1.66	0.77
1:A:94:ASN:HD21	1:A:217:THR:HA	1.49	0.77
1:B:201:MIS:H21	1:B:219:SER:HB2	1.68	0.76
1:A:75:GLU:HG3	1:A:76:ILE:H	1.51	0.74
1:B:65:GLY:HA2	1:B:102:ASP:HB2	1.69	0.73
1:B:261:ILE:HG12	1:B:284:VAL:HG11	1.69	0.73
1:C:258:TYR:CZ	1:C:260:GLY:HA2	2.24	0.73
1:A:305:ILE:CD1	1:A:336:MET:HG2	2.18	0.73
1:B:287:VAL:HG13	1:B:293:ALA:HB3	1.71	0.72
1:B:284:VAL:O	1:B:300:VAL:O	2.06	0.72
1:A:65:GLY:HA3	1:A:77:ARG:HE	1.53	0.71
1:C:265:GLU:OE1	1:C:314:SER:HA	1.90	0.71
1:A:245:MET:HG2	1:A:249:ILE:HD11	1.72	0.71
1:A:207:ASN:HB3	1:A:213:MET:HE2	1.72	0.70
1:A:170:GLN:HB2	1:B:172:ILE:HD13	1.72	0.70
1:B:179:ILE:HA	1:B:187:GLN:O	1.93	0.69
1:A:240:LEU:O	1:A:244:ILE:HG12	1.93	0.69
1:C:330:VAL:HG22	1:C:347:THR:HG22	1.73	0.69
1:A:245:MET:O	1:A:249:ILE:HG13	1.93	0.68
1:C:213:MET:HE2	1:C:213:MET:HA	1.74	0.68
1:B:167:THR:HG23	1:C:174:SER:HB3	1.76	0.68
1:B:47:TYR:HB3	1:B:156:LEU:HD11	1.76	0.68
1:C:196:ILE:HG23	1:C:201:MIS:H31	1.75	0.67
1:A:261:ILE:HG23	1:A:284:VAL:HG13	1.76	0.67
1:B:92:ILE:HD11	1:B:245:MET:HB2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:ALA:CB	1:C:346:VAL:HG12	2.26	0.66
1:C:302:ASP:HB3	1:C:337:ARG:HG3	1.78	0.66
1:A:240:LEU:CD1	1:A:244:ILE:HD11	2.27	0.65
1:B:187:GLN:OE1	1:B:187:GLN:HA	1.95	0.65
1:C:281:GLY:C	1:C:303:LEU:HD11	2.22	0.65
1:A:305:ILE:HD11	1:A:336:MET:CG	2.23	0.65
1:B:124:LEU:HD22	1:B:256:ARG:HH21	1.61	0.64
1:B:64:ARG:NH2	1:B:74:LEU:HD13	2.09	0.64
1:B:308:ASP:HB2	1:B:331:ILE:HD13	1.77	0.64
1:A:258:TYR:CZ	1:A:260:GLY:HA2	2.32	0.64
1:A:351:TYR:CD2	1:A:352:PRO:HD2	2.33	0.63
1:C:258:TYR:O	1:C:292:PRO:HG2	1.99	0.63
1:C:314:SER:OG	1:C:317:GLU:HB3	1.98	0.63
1:A:201:MIS:H21	1:A:219:SER:HB2	1.81	0.62
1:A:240:LEU:HD12	1:A:244:ILE:HD11	1.81	0.62
1:B:95:LYS:HG3	1:B:121:SER:HB2	1.81	0.62
1:A:190:LEU:HG	1:A:237:PRO:HG3	1.81	0.62
1:A:308:ASP:HA	1:A:332:PRO:HD2	1.82	0.61
1:C:302:ASP:HB3	1:C:337:ARG:CG	2.30	0.61
1:C:259:ILE:HD11	2:F:410:PHE:CE2	2.34	0.61
1:B:287:VAL:HG21	1:B:298:ILE:HG23	1.82	0.61
1:A:50:ALA:HB2	1:A:208:SER:O	2.01	0.61
1:B:261:ILE:HD11	1:B:298:ILE:HD13	1.82	0.61
1:A:94:ASN:HD22	1:A:201:MIS:HB2	1.66	0.61
1:A:94:ASN:ND2	1:A:217:THR:HA	2.16	0.61
1:A:196:ILE:CG2	1:A:201:MIS:H31	2.31	0.61
1:A:64:ARG:NH2	1:A:66:LEU:HD11	2.16	0.60
1:C:284:VAL:HG12	1:C:300:VAL:HA	1.83	0.60
1:A:167:THR:HG23	1:B:174:SER:HB3	1.83	0.60
1:A:99:ASN:HD22	1:A:100:ASP:H	1.49	0.59
1:B:325:ILE:HD13	1:B:331:ILE:HD12	1.85	0.59
1:A:253:ARG:NH1	1:A:255:ILE:HG12	2.18	0.59
1:C:261:ILE:HG12	1:C:293:ALA:HB2	1.84	0.59
1:A:65:GLY:CA	1:A:77:ARG:HE	2.16	0.59
1:B:287:VAL:HG13	1:B:293:ALA:CB	2.33	0.59
1:C:261:ILE:CG1	1:C:293:ALA:HB2	2.33	0.58
1:B:261:ILE:CD1	1:B:298:ILE:HD13	2.32	0.58
1:A:176:THR:HG22	1:A:190:LEU:HD22	1.86	0.58
1:B:261:ILE:HG23	1:B:284:VAL:HG13	1.86	0.58
1:B:175:ALA:CB	1:B:178:ARG:HE	2.17	0.58
1:C:292:PRO:HG3	1:C:349:GLN:HG3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:ILE:HG12	1:A:168:ILE:HD12	1.84	0.58
1:C:223:SER:HG	1:C:227:GLU:N	2.02	0.58
1:B:79:LEU:CD2	1:B:163:ASN:HB2	2.34	0.57
1:A:330:VAL:HG22	1:A:347:THR:HG22	1.87	0.57
1:A:201:MIS:C2	1:A:219:SER:HB2	2.34	0.56
1:C:129:VAL:HG23	1:C:248:LEU:HD12	1.86	0.56
1:A:325:ILE:HD13	1:A:331:ILE:HG21	1.88	0.56
1:B:316:LEU:HD12	1:B:317:GLU:N	2.20	0.56
1:A:90:TYR:CE2	1:A:119:VAL:HG21	2.40	0.56
1:A:187:GLN:HB3	1:A:189:PHE:CE2	2.41	0.56
1:A:283:VAL:HG22	1:A:284:VAL:H	1.71	0.56
1:B:189:PHE:HD1	1:B:235:ALA:O	1.89	0.56
1:A:99:ASN:HD22	1:A:100:ASP:N	2.04	0.55
1:C:283:VAL:HG22	1:C:303:LEU:HD13	1.87	0.55
1:A:283:VAL:HG22	1:A:284:VAL:N	2.21	0.55
1:C:284:VAL:O	1:C:300:VAL:O	2.24	0.55
1:A:84:ILE:HD13	1:A:91:ILE:HG23	1.89	0.55
1:A:239:GLN:CD	1:A:239:GLN:H	2.13	0.55
1:B:287:VAL:HG11	1:B:294:ALA:HA	1.89	0.55
1:A:134:ALA:HB2	1:A:138:LEU:HD11	1.89	0.55
1:C:116:ALA:HB1	1:C:130:LEU:HB3	1.88	0.55
1:A:47:TYR:CZ	1:A:154:VAL:HG11	2.42	0.54
1:B:84:ILE:HG21	1:B:138:LEU:HD13	1.88	0.54
1:C:67:ASN:O	1:C:68:THR:C	2.51	0.54
1:C:261:ILE:HG12	1:C:293:ALA:CB	2.37	0.54
1:C:281:GLY:O	1:C:303:LEU:HD11	2.06	0.54
1:B:322:VAL:HG11	2:E:410:PHE:CD1	2.43	0.54
1:C:63:ASN:C	1:C:63:ASN:OD1	2.51	0.54
1:B:295:ASN:C	1:B:297:GLY:H	2.17	0.53
1:B:67:ASN:HB2	1:B:75:GLU:OE2	2.09	0.53
1:A:45:ALA:H	1:B:208:SER:HB3	1.73	0.53
1:A:209:LEU:HD11	1:C:44:PRO:HD3	1.91	0.53
1:A:141:ILE:HD12	1:A:142:PRO:HD2	1.90	0.53
1:A:251:ASP:N	1:A:251:ASP:OD1	2.41	0.53
1:C:262:GLY:HA2	2:F:408:TYR:O	2.08	0.53
1:A:250:ARG:C	1:A:251:ASP:OD1	2.52	0.53
1:C:92:ILE:HD11	1:C:245:MET:HB2	1.91	0.52
1:B:209:LEU:HB2	1:B:211:GLU:HG3	1.90	0.52
1:C:259:ILE:HG13	1:C:261:ILE:HG13	1.91	0.52
1:B:141:ILE:O	1:B:143:ILE:HG13	2.10	0.52
1:A:258:TYR:OH	1:A:260:GLY:HA2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:TYR:HE2	1:A:119:VAL:HG21	1.73	0.52
1:B:175:ALA:HB3	1:B:191:GLN:HB3	1.90	0.52
1:A:85:MET:HG3	1:A:92:ILE:HD11	1.90	0.52
1:A:98:ILE:HD12	1:A:98:ILE:C	2.35	0.52
1:C:292:PRO:HB3	1:C:347:THR:O	2.10	0.52
1:A:305:ILE:HD11	1:A:336:MET:N	2.25	0.52
1:A:179:ILE:HD12	1:A:183:PRO:CA	2.34	0.51
1:B:168:ILE:O	1:C:174:SER:HA	2.09	0.51
1:A:47:TYR:HB3	1:A:156:LEU:HD11	1.92	0.51
1:B:94:ASN:HB2	1:B:97:VAL:HG23	1.92	0.51
1:A:248:LEU:O	1:A:251:ASP:O	2.29	0.51
1:B:79:LEU:HD21	1:B:163:ASN:HB2	1.92	0.51
1:A:137:GLY:O	1:A:138:LEU:HD23	2.09	0.51
1:A:293:ALA:HB1	1:A:298:ILE:HD12	1.92	0.51
1:A:55:ALA:HB1	1:A:166:GLN:HE22	1.76	0.51
1:A:328:GLY:N	1:A:348:ILE:O	2.40	0.51
1:B:344:LEU:HD12	1:B:344:LEU:O	2.11	0.51
1:B:213:MET:HA	1:B:213:MET:HE2	1.94	0.50
1:A:194:ALA:O	1:A:196:ILE:HG12	2.11	0.50
1:B:310:LYS:HG3	1:B:321:GLN:OE1	2.12	0.50
1:B:122:ASP:CG	1:B:254:VAL:HG11	2.36	0.50
1:B:299:GLN:HG2	1:B:300:VAL:H	1.77	0.50
1:B:322:VAL:HG11	2:E:410:PHE:HD1	1.77	0.50
1:C:305:ILE:HD11	1:C:336:MET:CG	2.41	0.50
1:A:47:TYR:CE2	1:A:154:VAL:HG11	2.47	0.50
1:C:316:LEU:HG	1:C:319:MET:HE2	1.93	0.50
1:A:65:GLY:HA3	1:A:77:ARG:NE	2.26	0.49
1:A:166:GLN:O	1:B:178:ARG:NH2	2.46	0.49
1:C:47:TYR:HB3	1:C:156:LEU:HD11	1.94	0.49
1:A:65:GLY:C	1:A:77:ARG:HE	2.21	0.49
1:A:85:MET:HE1	1:A:212:LEU:CD2	2.42	0.49
1:A:86:ASP:OD1	1:A:90:TYR:HB2	2.13	0.49
1:A:189:PHE:HB3	1:A:234:PHE:HB3	1.94	0.49
1:A:251:ASP:C	1:A:253:ARG:H	2.21	0.49
1:C:85:MET:HG3	1:C:245:MET:SD	2.53	0.49
1:C:178:ARG:O	1:C:189:PHE:N	2.45	0.49
1:B:258:TYR:CZ	1:B:260:GLY:HA2	2.47	0.49
1:A:103:GLN:NE2	1:A:105:ILE:HD11	2.27	0.49
1:A:304:ILE:HG12	1:A:335:VAL:HG12	1.93	0.49
1:B:178:ARG:HG3	1:B:191:GLN:HB2	1.94	0.49
1:A:237:PRO:HB2	1:A:239:GLN:OE1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:ASN:HB2	1:C:97:VAL:HG23	1.95	0.48
1:A:230:GLU:OE2	1:C:231:GLY:N	2.42	0.48
1:A:64:ARG:HB3	1:A:64:ARG:NH1	2.29	0.48
1:A:209:LEU:HD22	1:C:42:GLU:HG2	1.96	0.48
1:A:255:ILE:HG23	1:A:350:GLU:HB2	1.96	0.48
1:B:223:SER:OG	1:B:227:GLU:N	2.47	0.48
1:B:251:ASP:C	1:B:253:ARG:H	2.21	0.48
1:B:259:ILE:HG23	1:B:261:ILE:H	1.78	0.48
1:A:64:ARG:NH1	1:A:103:GLN:OE1	2.46	0.48
1:A:90:TYR:HA	1:A:130:LEU:O	2.14	0.48
1:A:141:ILE:O	1:A:141:ILE:HG23	2.14	0.48
1:A:292:PRO:HG3	1:A:349:GLN:CG	2.30	0.48
1:C:305:ILE:HD11	1:C:336:MET:SD	2.54	0.48
1:C:58:VAL:HG11	1:C:158:ILE:HG22	1.96	0.47
1:A:85:MET:HE1	1:A:212:LEU:HD22	1.96	0.47
1:C:90:TYR:CE2	1:C:131:LYS:HD3	2.49	0.47
1:C:206:VAL:HG12	1:C:212:LEU:HA	1.95	0.47
1:A:134:ALA:HB1	1:A:138:LEU:HD21	1.95	0.47
1:A:95:LYS:HD3	1:A:123:SER:HA	1.96	0.47
1:B:84:ILE:HG12	1:B:91:ILE:HG12	1.97	0.47
1:B:260:GLY:HA3	1:B:288:SER:HB3	1.97	0.47
1:A:248:LEU:HA	1:A:251:ASP:O	2.14	0.47
1:A:305:ILE:HD12	1:A:305:ILE:N	2.29	0.47
1:A:207:ASN:OD1	1:A:211:GLU:N	2.48	0.47
1:B:177:GLY:HA2	1:B:188:ASN:CG	2.40	0.47
1:A:94:ASN:OD1	1:A:217:THR:HG23	2.15	0.47
1:A:147:ARG:HG2	1:A:211:GLU:OE1	2.14	0.47
1:B:158:ILE:HG12	1:B:168:ILE:HD12	1.97	0.47
1:C:300:VAL:HG23	1:C:301:ASN:N	2.30	0.47
1:A:99:ASN:ND2	1:A:100:ASP:N	2.64	0.46
1:B:79:LEU:HD22	1:B:163:ASN:HB2	1.97	0.46
1:B:287:VAL:CG1	1:B:294:ALA:N	2.78	0.46
1:A:94:ASN:HB2	1:A:97:VAL:HG23	1.97	0.46
1:A:230:GLU:O	1:C:195:SER:HB3	2.16	0.46
1:A:220:PHE:HD2	1:A:229:PRO:HG2	1.79	0.46
1:A:159:GLY:C	1:A:161:PRO:HD3	2.41	0.46
1:A:300:VAL:O	1:A:301:ASN:HB2	2.15	0.46
1:B:299:GLN:HG2	1:B:300:VAL:N	2.30	0.46
1:C:284:VAL:CG1	1:C:300:VAL:HA	2.44	0.46
1:A:320:ALA:O	1:A:324:GLU:HG2	2.15	0.46
1:B:179:ILE:O	1:B:179:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:GLN:OE1	1:B:187:GLN:CA	2.60	0.46
1:C:304:ILE:HG12	1:C:335:VAL:HG12	1.98	0.46
1:A:263:GLY:HA3	1:A:283:VAL:O	2.16	0.46
1:B:261:ILE:HG13	1:B:293:ALA:CB	2.45	0.46
1:C:196:ILE:CG2	1:C:201:MIS:H31	2.46	0.46
1:B:124:LEU:O	1:B:187:GLN:NE2	2.50	0.45
1:C:158:ILE:HG12	1:C:168:ILE:HD12	1.98	0.45
1:C:73:GLN:HG2	1:C:74:LEU:H	1.82	0.45
1:A:90:TYR:CZ	1:A:249:ILE:HG23	2.50	0.45
1:B:254:VAL:HG13	1:B:254:VAL:O	2.16	0.45
1:B:295:ASN:O	1:B:296:ALA:HB3	2.17	0.45
1:A:65:GLY:C	1:A:66:LEU:HD12	2.41	0.45
1:B:319:MET:HA	1:B:322:VAL:HG12	1.99	0.45
1:B:322:VAL:HG23	1:B:325:ILE:HD12	1.98	0.45
1:A:114:PHE:CD1	1:A:132:ILE:HD12	2.51	0.45
1:A:165:GLY:HA3	1:B:178:ARG:HH12	1.82	0.45
1:A:207:ASN:HB3	1:A:213:MET:CE	2.41	0.45
1:B:66:LEU:HD23	1:B:74:LEU:HG	1.99	0.45
1:A:83:VAL:CG1	1:A:141:ILE:HB	2.47	0.44
1:A:220:PHE:CE2	1:A:222:LYS:CB	3.01	0.44
1:B:177:GLY:HA2	1:B:188:ASN:OD1	2.17	0.44
1:C:73:GLN:HG2	1:C:74:LEU:N	2.32	0.44
1:A:94:ASN:CG	1:A:217:THR:HG23	2.42	0.44
1:B:285:ASN:OD1	1:B:286:GLU:N	2.51	0.44
1:B:287:VAL:HG11	1:B:294:ALA:CA	2.47	0.44
1:B:288:SER:O	1:B:291:GLY:N	2.48	0.44
1:C:73:GLN:CG	1:C:74:LEU:H	2.30	0.44
1:B:178:ARG:HD3	1:B:178:ARG:HA	1.87	0.44
1:C:214:GLY:HA2	1:C:236:ILE:O	2.18	0.44
1:A:42:GLU:HB2	1:B:209:LEU:HD22	2.00	0.44
1:A:95:LYS:HE3	1:A:123:SER:OG	2.18	0.43
1:A:305:ILE:HD12	1:A:335:VAL:HA	1.98	0.43
1:B:250:ARG:HG2	1:B:251:ASP:OD1	2.18	0.43
1:C:314:SER:OG	1:C:317:GLU:CB	2.66	0.43
1:B:67:ASN:ND2	1:B:75:GLU:OE2	2.43	0.43
1:B:310:LYS:HA	1:B:311:PRO:HD3	1.88	0.43
1:C:142:PRO:HB2	1:C:211:GLU:HA	2.00	0.43
1:A:351:TYR:OH	2:D:409:ARG:O	2.26	0.43
1:A:174:SER:HA	1:C:168:ILE:O	2.17	0.43
1:A:225:ASP:OD1	1:A:225:ASP:N	2.52	0.43
1:C:254:VAL:HG11	1:C:256:ARG:CZ	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:ALA:CB	1:B:178:ARG:NE	2.81	0.43
1:A:94:ASN:HA	1:A:126:ASP:O	2.19	0.43
1:B:178:ARG:C	1:B:179:ILE:HG23	2.44	0.43
1:C:124:LEU:HD12	1:C:124:LEU:HA	1.85	0.42
1:C:263:GLY:HA3	1:C:283:VAL:O	2.18	0.42
1:A:240:LEU:HD11	1:A:244:ILE:HD11	1.98	0.42
1:A:250:ARG:CZ	1:A:251:ASP:OD1	2.66	0.42
1:A:63:ASN:OD1	1:A:63:ASN:C	2.63	0.42
1:C:60:ASN:HB2	1:C:107:ALA:HB3	2.00	0.42
1:C:258:TYR:CZ	1:C:260:GLY:CA	3.01	0.42
1:A:38:ASP:HB3	1:A:40:THR:HG22	2.01	0.42
1:A:55:ALA:HB1	1:A:166:GLN:NE2	2.34	0.42
1:A:55:ALA:HB3	1:A:56:PRO:HD3	2.01	0.42
1:A:43:THR:HA	1:A:44:PRO:HD3	1.83	0.42
1:B:155:VAL:HA	1:B:206:VAL:O	2.20	0.42
1:A:352:PRO:O	1:A:353:ALA:C	2.63	0.42
1:B:48:ASN:O	1:B:52:ARG:HG3	2.20	0.42
1:B:50:ALA:HB2	1:B:208:SER:O	2.20	0.42
1:C:305:ILE:HD11	1:C:336:MET:HG2	2.01	0.42
1:B:223:SER:HB2	1:B:227:GLU:O	2.20	0.42
1:A:197:ASN:HB3	1:A:198:PRO:HD2	2.02	0.42
1:B:259:ILE:HG23	1:B:261:ILE:N	2.35	0.42
1:C:58:VAL:HB	1:C:166:GLN:NE2	2.35	0.42
1:C:282:ILE:HD13	1:C:315:ALA:HA	2.02	0.42
1:A:178:ARG:HH22	1:C:166:GLN:C	2.28	0.42
1:A:187:GLN:HB3	1:A:189:PHE:CD2	2.55	0.42
1:A:64:ARG:HB3	1:A:64:ARG:CZ	2.50	0.41
1:A:120:GLY:HA3	1:A:248:LEU:HD22	2.02	0.41
1:B:47:TYR:CE2	1:B:154:VAL:HG11	2.55	0.41
1:B:63:ASN:HB2	1:B:97:VAL:O	2.19	0.41
1:B:298:ILE:HG23	1:B:298:ILE:O	2.19	0.41
1:C:61:VAL:HG22	1:C:106:VAL:HG22	2.01	0.41
1:A:56:PRO:HB2	1:A:139:PRO:HB3	2.01	0.41
1:A:255:ILE:CG2	1:A:327:PRO:HG2	2.50	0.41
1:B:161:PRO:HA	1:B:198:PRO:O	2.20	0.41
1:C:197:ASN:HB3	1:C:198:PRO:HD2	2.01	0.41
1:C:218:LEU:HB2	1:C:234:PHE:HB2	2.01	0.41
1:C:286:GLU:HG3	2:F:409:ARG:NH2	2.35	0.41
1:C:85:MET:HG3	1:C:92:ILE:HG13	2.02	0.41
1:C:330:VAL:HA	1:C:346:VAL:O	2.19	0.41
1:A:148:VAL:O	1:A:150:HIS:ND1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:PHE:C	1:A:220:PHE:CD2	2.98	0.41
1:C:98:ILE:C	1:C:98:ILE:HD12	2.46	0.41
1:A:84:ILE:O	1:A:141:ILE:HG22	2.21	0.41
1:A:144:ASN:OD1	1:A:144:ASN:C	2.64	0.41
1:C:106:VAL:HG23	1:C:116:ALA:HB2	2.03	0.41
1:C:282:ILE:O	1:C:303:LEU:HD12	2.20	0.41
1:B:116:ALA:HB1	1:B:130:LEU:HB3	2.03	0.41
1:B:174:SER:HB2	1:B:191:GLN:HG2	2.02	0.41
1:B:261:ILE:HG12	1:B:284:VAL:CG1	2.44	0.40
1:C:63:ASN:ND2	1:C:101:ALA:HB2	2.35	0.40
1:A:161:PRO:HG3	1:A:200:ASN:OD1	2.21	0.40
1:A:172:ILE:HD13	1:C:170:GLN:HB2	2.02	0.40
1:A:245:MET:HG2	1:A:249:ILE:CD1	2.48	0.40
1:B:223:SER:CB	1:B:227:GLU:O	2.70	0.40
1:C:251:ASP:C	1:C:253:ARG:H	2.29	0.40
1:C:308:ASP:HA	1:C:332:PRO:HD2	2.04	0.40
1:A:93:THR:O	1:A:128:ALA:N	2.55	0.40
1:B:184:THR:OG1	1:B:185:GLY:N	2.53	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	288/340 (85%)	274 (95%)	14 (5%)	0	100	100
1	B	266/340 (78%)	253 (95%)	13 (5%)	0	100	100
1	C	282/340 (83%)	277 (98%)	5 (2%)	0	100	100
2	D	1/3 (33%)	0	1 (100%)	0	100	100
2	E	1/3 (33%)	0	1 (100%)	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	839/1029 (82%)	805 (96%)	34 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	232/275 (84%)	224 (97%)	8 (3%)	32	66
1	B	223/275 (81%)	221 (99%)	2 (1%)	70	85
1	C	235/275 (86%)	234 (100%)	1 (0%)	84	90
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	699/834 (84%)	688 (98%)	11 (2%)	55	79

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	THR
1	A	91	ILE
1	A	93	THR
1	A	206	VAL
1	A	225	ASP
1	A	239	GLN
1	A	251	ASP
1	A	349	GLN
1	B	261	ILE
1	B	346	VAL
1	C	196	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	ASN
1	A	99	ASN
1	A	109	GLN
1	A	166	GLN
1	A	295	ASN
1	A	321	GLN
1	A	345	GLN
1	B	166	GLN
1	C	87	GLN
1	C	295	ASN

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	B	201	1	11,12,13	0.69	0	11,16,18	0.85	0
1	MIS	A	201	1	11,12,13	0.68	0	11,16,18	0.62	0
1	MIS	C	201	1	11,12,13	0.69	0	11,16,18	0.60	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	B	201	1	-	5/11/13/15	-
1	MIS	A	201	1	-	4/11/13/15	-
1	MIS	C	201	1	-	5/11/13/15	-

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	C1-O3P-P-O1P
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-O2P
1	C	201	MIS	CB-OG-P-O3P
1	C	201	MIS	C1-O3P-P-O1P
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	A	201	MIS	C1-O3P-P-O2P
1	C	201	MIS	C1-O3P-P-O2P

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	201	MIS	1	0
1	A	201	MIS	4	0
1	C	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 ⓘ

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	294/340 (86%)	0.01	3 (1%) 79 59	66, 133, 179, 197	0
1	B	278/340 (81%)	0.03	4 (1%) 73 51	51, 81, 206, 234	0
1	C	292/340 (85%)	-0.06	0 100 100	60, 100, 151, 184	0
2	D	3/3 (100%)	0.35	0 100 100	175, 175, 181, 199	0
2	E	3/3 (100%)	0.51	0 100 100	205, 205, 212, 219	0
2	F	3/3 (100%)	0.06	0 100 100	117, 117, 152, 155	0
All	All	873/1029 (84%)	-0.01	7 (0%) 82 64	51, 105, 181, 234	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	213	MET	2.9
1	A	353	ALA	2.6
1	A	85	MET	2.2
1	B	352	PRO	2.2
1	A	185	GLY	2.2
1	B	297	GLY	2.1
1	B	224	ASN	2.1

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	0.95	0.09	71,86,98,105	0
1	MIS	C	201	13/14	0.97	0.07	57,64,78,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	MIS	B	201	13/14	0.98	0.07	48,49,66,66	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.