



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 05:54 AM UTC

PDB ID : 2Z9J / pdb_00002z9j
Title : Complex structure of SARS-CoV 3C-like protease with EPDTC
Authors : Lee, C.C.; Wang, A.H.
Deposited on : 2007-09-20
Resolution : 1.95 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

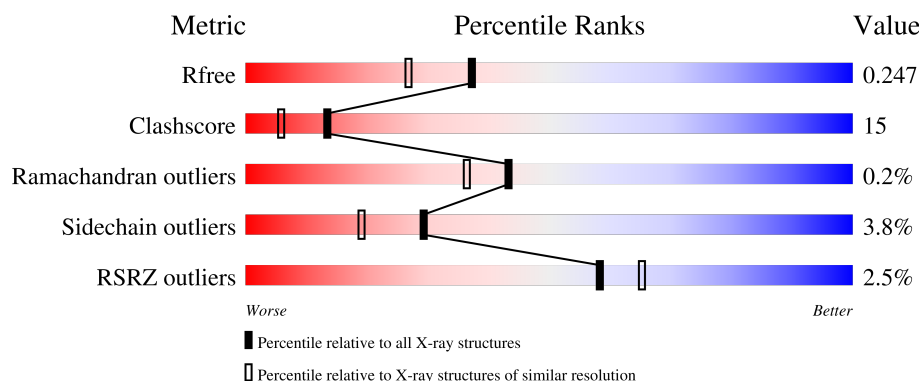
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	306	
1	B	306	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DMS	A	802	-	X	-	-
2	DMS	A	804	-	X	-	-
2	DMS	B	801	-	X	X	-
2	DMS	B	803	-	X	-	-

2 Entry composition [i](#)

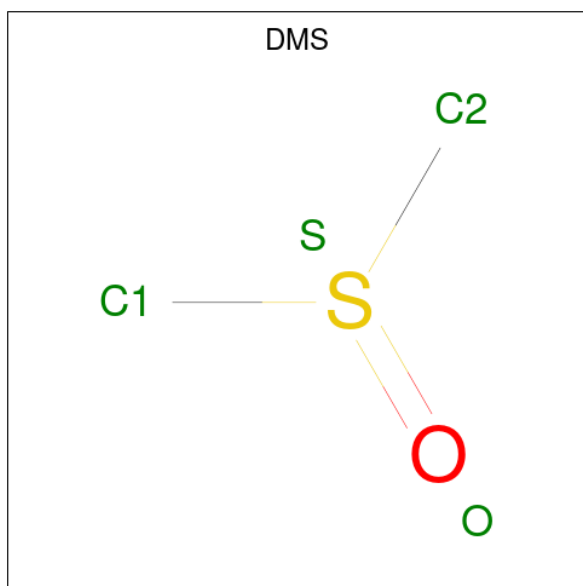
There are 4 unique types of molecules in this entry. The entry contains 5440 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3C-like proteinase.

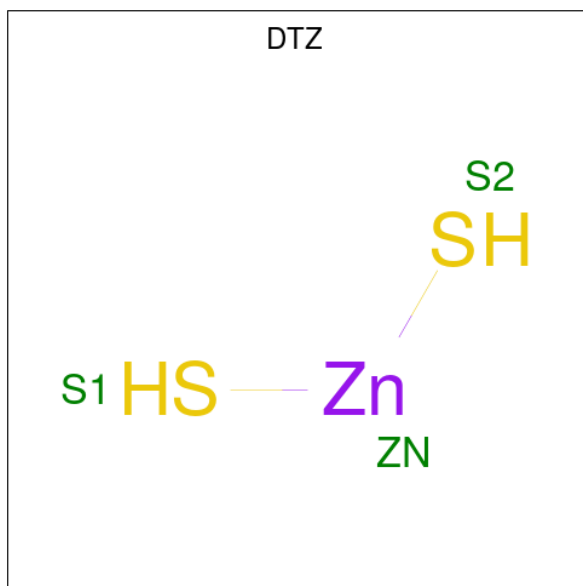
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2336	1476	400	438	22			
1	B	304	Total	C	N	O	S	0	0	0
			2350	1485	402	441	22			

- Molecule 2 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

- Molecule 3 is zinc(II)hydrosulfide (CCD ID: DTZ) (formula: $\text{H}_2\text{S}_2\text{Zn}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	S	Zn	0	0
			3	2	1		
3	B	1	Total	S	Zn	0	0
			3	2	1		

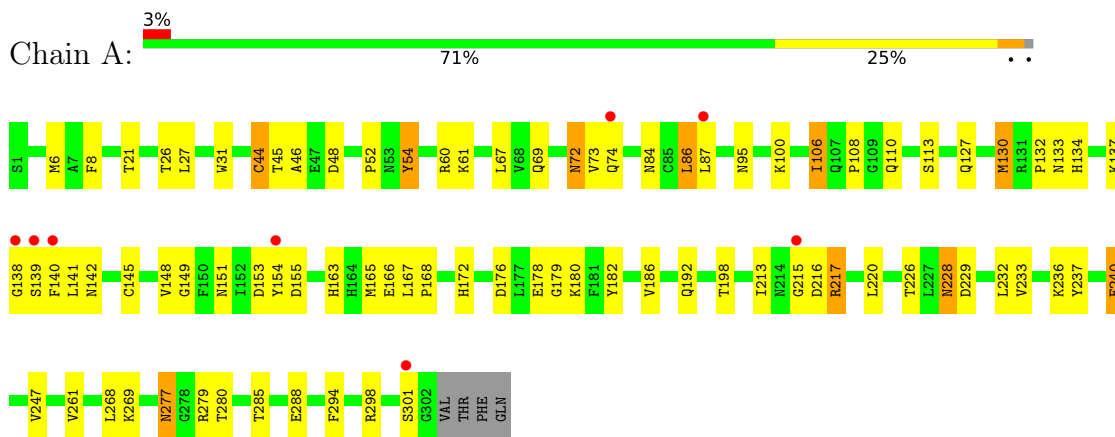
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	381	Total	O	0	0
			381	381		
4	B	351	Total	O	0	0
			351	351		

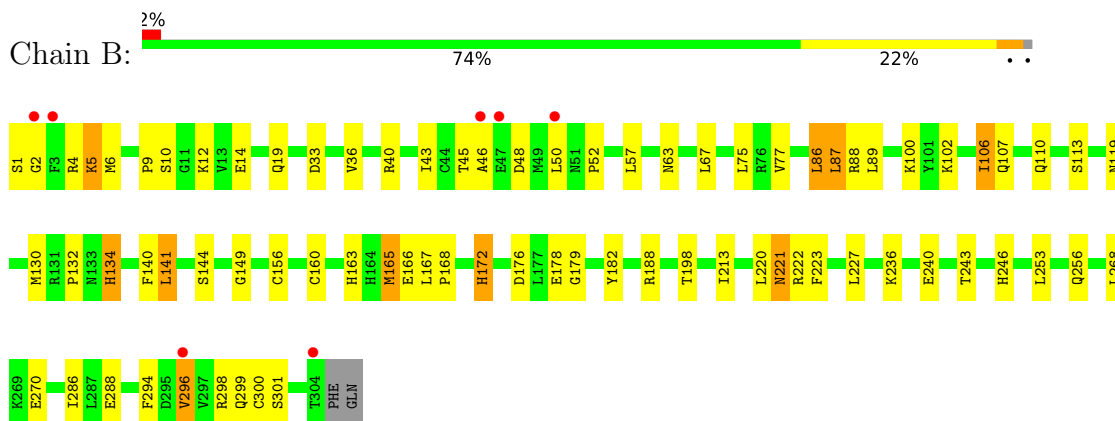
3 Residue-property plots

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: 3C-like proteinase



• Molecule 1: 3C-like proteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.15Å 96.30Å 67.62Å 90.00° 103.63° 90.00°	Depositor
Resolution (Å)	30.00 – 1.95 30.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.95) 93.6 (30.00-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.194 , 0.248 0.191 , 0.247	Depositor DCC
R_{free} test set	2361 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtriage
Anisotropy	0.393	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5440	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: DTZ, DMS

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	0.51	0/2388	0.99	12/3244 (0.4%)
1	B	0.51	0/2402	0.98	10/3264 (0.3%)
All	All	0.51	0/4790	0.98	22/6508 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	LYS	N-CA-C	-7.41	98.84	109.96
1	A	178	GLU	N-CA-C	-7.20	103.66	113.30
1	B	176	ASP	N-CA-C	-6.76	101.02	110.35
1	B	165	MET	N-CA-C	6.55	118.16	108.60
1	A	182	TYR	N-CA-C	-6.39	98.86	109.46
1	A	130	MET	N-CA-C	-6.31	99.59	109.25
1	B	182	TYR	N-CA-C	-6.29	99.02	109.46
1	A	138	GLY	N-CA-C	6.24	127.97	113.18
1	A	73	VAL	N-CA-C	6.24	117.28	108.36
1	B	100	LYS	N-CA-C	-6.11	101.24	110.28
1	B	134	HIS	N-CA-C	6.06	120.62	113.23
1	A	217	ARG	N-CA-C	5.94	119.76	111.56
1	A	100	LYS	N-CA-C	-5.88	101.15	109.96
1	B	130	MET	N-CA-C	-5.84	101.00	109.59
1	B	178	GLU	N-CA-C	-5.64	105.74	113.30
1	B	172	HIS	N-CA-C	5.44	118.43	109.72
1	A	165	MET	N-CA-C	5.44	116.54	108.60
1	A	44	CYS	N-CA-C	5.43	118.77	110.20
1	A	54	TYR	N-CA-C	5.40	117.17	111.28
1	A	148	VAL	N-CA-C	5.37	117.36	108.99
1	B	9	PRO	N-CA-C	-5.33	102.82	111.14
1	A	176	ASP	N-CA-C	-5.22	102.27	110.36

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	2336	0	2288	74	0
1	B	2350	0	2304	69	0
2	A	8	0	12	2	0
2	B	8	0	12	6	0
3	A	3	0	0	1	0
3	B	3	0	0	0	0
4	A	381	0	0	18	0
4	B	351	0	0	9	0
All	All	5440	0	4616	137	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 15.

All (137) 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:139:SER:HA	1:B:2:GLY:H	1.16	1.08
1:B:163:HIS:NE2	2:B:801:DMS:H22	1.72	1.04
1:B:45:THR:HG22	1:B:46:ALA:H	1.26	1.01
1:B:36:VAL:CG1	1:B:89:LEU:HB2	2.07	0.85
1:B:221:ASN:C	1:B:221:ASN:HD22	1.87	0.82
1:B:45:THR:HB	4:B:1228:HOH:O	1.80	0.81
1:B:36:VAL:HG13	1:B:89:LEU:HB2	1.62	0.80
1:B:77:VAL:HG12	4:B:1028:HOH:O	1.82	0.79
1:A:139:SER:HA	1:B:2:GLY:N	1.97	0.77
1:A:139:SER:HB3	4:B:1061:HOH:O	1.85	0.77
1:A:21:THR:HG23	4:A:1140:HOH:O	1.87	0.74
1:A:226:THR:HG23	1:A:229:ASP:H	1.52	0.72
1:B:270:GLU:HG3	4:B:1063:HOH:O	1.89	0.71
1:A:145:CYS:SG	2:A:804:DMS:H22	2.33	0.68
1:A:21:THR:HG22	1:A:26:THR:OG1	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ASP:HB3	4:A:1189:HOH:O	1.93	0.68
1:A:198:THR:OG1	1:A:240:GLU:HG2	1.93	0.68
1:A:236:LYS:O	1:A:236:LYS:HD3	1.94	0.68
1:B:45:THR:HG22	1:B:46:ALA:N	2.04	0.68
1:A:45:THR:HG22	1:A:46:ALA:H	1.59	0.68
1:A:6:MET:HE1	4:B:1026:HOH:O	1.95	0.67
1:A:67:LEU:HD11	1:A:74:GLN:NE2	2.10	0.67
1:A:21:THR:OG1	1:A:67:LEU:HB3	1.94	0.66
1:A:186:VAL:H	1:A:192:GLN:HE22	1.41	0.66
1:A:279:ARG:HD2	4:A:1247:HOH:O	1.96	0.65
1:A:69:GLN:HG2	1:A:74:GLN:NE2	2.12	0.65
1:A:72:ASN:HA	4:A:1141:HOH:O	1.98	0.63
1:A:74:GLN:HG2	4:A:1256:HOH:O	1.97	0.63
1:A:226:THR:CG2	1:A:229:ASP:H	2.12	0.62
1:A:84:ASN:CG	1:A:180:LYS:HD3	2.25	0.62
4:A:1235:HOH:O	1:B:4:ARG:HD3	2.01	0.61
1:B:166:GLU:HB2	2:B:801:DMS:C2	2.30	0.60
1:A:134:HIS:HD2	4:A:1151:HOH:O	1.84	0.59
1:B:52:PRO:HB2	1:B:57:LEU:HD11	1.83	0.59
1:B:50:LEU:HD12	1:B:50:LEU:N	2.18	0.59
1:B:166:GLU:HB2	2:B:801:DMS:H23	1.86	0.58
1:B:45:THR:CG2	1:B:46:ALA:H	2.09	0.57
1:B:33:ASP:OD2	4:B:1003:HOH:O	2.17	0.57
1:B:10:SER:O	1:B:14:GLU:HG3	2.05	0.56
1:B:294:PHE:HD2	1:B:298:ARG:NH2	2.04	0.56
1:A:27:LEU:HD23	3:A:902:DTZ:S1	2.46	0.56
1:A:294:PHE:CD2	1:A:298:ARG:NH2	2.74	0.56
1:B:52:PRO:HB2	1:B:57:LEU:CD1	2.35	0.56
1:A:86:LEU:HD22	1:A:179:GLY:CA	2.36	0.56
1:A:240:GLU:HG3	4:A:1259:HOH:O	2.06	0.56
1:A:236:LYS:HD2	1:A:237:TYR:CE1	2.42	0.55
1:B:106:ILE:HD13	1:B:160:CYS:CB	2.37	0.55
1:A:110:GLN:HG2	4:A:1245:HOH:O	2.06	0.54
1:A:137:LYS:HE3	4:A:1222:HOH:O	2.08	0.54
1:B:134:HIS:HD2	4:B:1000:HOH:O	1.89	0.54
1:A:6:MET:HE1	4:A:1264:HOH:O	2.06	0.54
1:B:141:LEU:C	2:B:801:DMS:H13	2.33	0.53
1:B:48:ASP:O	1:B:52:PRO:HB3	2.09	0.53
1:A:48:ASP:O	1:A:52:PRO:HB3	2.10	0.52
1:A:137:LYS:HE2	4:A:1232:HOH:O	2.09	0.52
1:B:106:ILE:HG13	1:B:110:GLN:HE21	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:ASN:ND2	1:B:223:PHE:H	2.07	0.52
1:A:167:LEU:HB3	1:A:168:PRO:HD2	1.92	0.52
1:B:198:THR:OG1	1:B:240:GLU:HG2	2.10	0.52
1:B:63:ASN:O	1:B:77:VAL:HG11	2.11	0.51
1:A:217:ARG:HB2	1:A:220:LEU:HD12	1.92	0.51
1:B:40:ARG:O	1:B:43:ILE:HG12	2.09	0.51
1:A:45:THR:HB	1:A:48:ASP:OD1	2.11	0.51
1:B:221:ASN:C	1:B:221:ASN:ND2	2.61	0.51
1:A:86:LEU:HD22	1:A:179:GLY:N	2.27	0.50
1:A:280:THR:HG22	1:A:285:THR:HG22	1.93	0.50
1:A:301:SER:HB2	4:A:1258:HOH:O	2.11	0.50
1:A:166:GLU:OE1	1:B:1:SER:N	2.43	0.50
1:A:236:LYS:HD3	1:A:236:LYS:C	2.36	0.50
1:A:61:LYS:HE2	4:A:940:HOH:O	2.12	0.50
1:B:107:GLN:H	1:B:110:GLN:NE2	2.10	0.49
1:A:228:ASN:O	1:A:232:LEU:HG	2.12	0.48
1:B:286:ILE:HD11	1:B:288:GLU:OE2	2.12	0.48
1:A:140:PHE:C	1:B:1:SER:HB2	2.37	0.48
1:B:166:GLU:N	2:B:801:DMS:H23	2.29	0.48
1:B:299:GLN:C	1:B:301:SER:H	2.21	0.48
1:B:19:GLN:HE21	1:B:119:ASN:HA	1.80	0.46
1:A:233:VAL:HG21	1:A:269:LYS:HD2	1.96	0.46
1:B:141:LEU:O	2:B:801:DMS:H13	2.15	0.46
1:A:31:TRP:CE2	1:A:95:ASN:HB2	2.51	0.46
1:A:142:ASN:ND2	2:A:804:DMS:H11	2.31	0.46
1:A:294:PHE:CE2	1:A:298:ARG:NH2	2.84	0.46
1:A:72:ASN:OD1	1:A:72:ASN:N	2.48	0.46
1:A:45:THR:HG22	1:A:46:ALA:N	2.28	0.45
1:B:165:MET:HE3	1:B:167:LEU:HD21	1.98	0.45
1:B:213:ILE:HD12	1:B:256:GLN:HE21	1.82	0.45
1:A:294:PHE:HD2	1:A:298:ARG:NH2	2.13	0.45
1:A:269:LYS:HE3	4:A:994:HOH:O	2.16	0.45
1:B:50:LEU:HD12	1:B:50:LEU:H	1.81	0.45
1:B:12:LYS:HE3	4:B:1241:HOH:O	2.16	0.44
1:A:45:THR:H	1:A:48:ASP:HB2	1.81	0.44
1:A:288:GLU:HB3	4:A:1016:HOH:O	2.17	0.44
1:A:8:PHE:HE2	1:A:151:ASN:HD22	1.65	0.44
1:B:294:PHE:HD2	1:B:298:ARG:HH21	1.66	0.44
1:A:163:HIS:CE1	1:A:172:HIS:HB3	2.52	0.44
1:B:106:ILE:HD13	1:B:160:CYS:HB3	2.00	0.44
1:A:247:VAL:HG13	1:A:261:VAL:HG11	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ARG:HD3	4:B:1176:HOH:O	2.18	0.44
1:B:213:ILE:HD12	1:B:256:GLN:NE2	2.33	0.44
1:B:86:LEU:HD13	1:B:179:GLY:HA2	1.99	0.43
1:B:243:THR:H	1:B:246:HIS:CD2	2.36	0.43
1:A:113:SER:O	1:A:149:GLY:HA2	2.18	0.43
1:B:163:HIS:HE1	1:B:172:HIS:HB3	1.83	0.43
1:B:50:LEU:N	1:B:50:LEU:CD1	2.82	0.43
1:A:216:ASP:OD2	1:A:216:ASP:C	2.61	0.43
1:A:45:THR:N	1:A:48:ASP:HB2	2.34	0.43
1:A:140:PHE:O	1:B:1:SER:HB2	2.19	0.42
1:A:213:ILE:C	1:A:215:GLY:H	2.26	0.42
1:B:163:HIS:CE1	1:B:172:HIS:HB3	2.54	0.42
1:B:222:ARG:O	1:B:222:ARG:HD3	2.19	0.42
1:A:141:LEU:HA	1:B:1:SER:HB2	2.00	0.42
1:A:277:ASN:HD22	1:A:277:ASN:HA	1.59	0.42
1:B:5:LYS:HB2	1:B:5:LYS:HE3	1.83	0.42
1:B:87:LEU:HA	1:B:87:LEU:HD22	1.79	0.42
1:A:21:THR:CG2	1:A:26:THR:OG1	2.63	0.42
1:A:6:MET:SD	4:A:1268:HOH:O	2.62	0.42
1:A:60:ARG:NH2	4:A:1070:HOH:O	2.52	0.41
1:A:108:PRO:HA	1:A:130:MET:CG	2.50	0.41
1:B:67:LEU:HD12	1:B:67:LEU:HA	1.92	0.41
1:A:153:ASP:O	1:A:154:TYR:HB2	2.20	0.41
1:B:86:LEU:HD22	1:B:179:GLY:N	2.35	0.41
1:B:113:SER:O	1:B:149:GLY:HA2	2.20	0.41
1:B:167:LEU:HB3	1:B:168:PRO:HD2	2.03	0.41
1:A:127:GLN:HE21	1:A:127:GLN:HA	1.84	0.41
1:B:6:MET:HE2	1:B:299:GLN:CD	2.45	0.41
1:B:102:LYS:HE3	1:B:156:CYS:SG	2.60	0.41
1:A:21:THR:HG22	1:A:26:THR:HG23	2.02	0.41
1:B:50:LEU:H	1:B:50:LEU:CD1	2.33	0.41
1:B:236:LYS:O	1:B:236:LYS:HD3	2.21	0.41
1:B:87:LEU:HD12	1:B:89:LEU:HD21	2.02	0.41
1:B:140:PHE:HB3	1:B:144:SER:OG	2.21	0.41
1:A:133:ASN:O	1:A:134:HIS:HB2	2.21	0.40
1:A:163:HIS:HE1	1:A:172:HIS:HB3	1.86	0.40
1:B:88:ARG:NH1	1:B:88:ARG:HB3	2.36	0.40
1:A:106:ILE:HG13	1:A:110:GLN:OE1	2.21	0.40
1:A:44:CYS:SG	1:A:54:TYR:CE2	3.15	0.40
1:B:253:LEU:HD11	1:B:296:VAL:HG21	2.04	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	300/306 (98%)	292 (97%)	8 (3%)	0	100	100
1	B	302/306 (99%)	289 (96%)	12 (4%)	1 (0%)	36	28
All	All	602/612 (98%)	581 (96%)	20 (3%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	300	CYS

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	259/263 (98%)	250 (96%)	9 (4%)	32	22
1	B	261/263 (99%)	250 (96%)	11 (4%)	26	16
All	All	520/526 (99%)	500 (96%)	20 (4%)	29	19

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	86	LEU
1	A	87	LEU
1	A	106	ILE
1	A	132	PRO

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Mol	Chain	Res	Type
1	A	228	ASN
1	A	240	GLU
1	A	268	LEU
1	A	277	ASN
1	B	75	LEU
1	B	86	LEU
1	B	87	LEU
1	B	106	ILE
1	B	132	PRO
1	B	141	LEU
1	B	220	LEU
1	B	221	ASN
1	B	227	LEU
1	B	268	LEU
1	B	296	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	95	ASN
1	A	119	ASN
1	A	127	GLN
1	A	134	HIS
1	A	142	ASN
1	A	151	ASN
1	A	192	GLN
1	A	256	GLN
1	A	277	ASN
1	B	19	GLN
1	B	51	ASN
1	B	83	GLN
1	B	110	GLN
1	B	134	HIS
1	B	189	GLN
1	B	221	ASN
1	B	256	GLN
1	B	273	GLN
1	B	277	ASN
1	B	299	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	DMS	A	804	-	3,3,3	0.27	0	3,3,3	3.27	3 (100%)
3	DTZ	B	901	1	0,2,2	-	-	-		
3	DTZ	A	902	1	0,2,2	-	-	-		
2	DMS	A	802	-	3,3,3	0.37	0	3,3,3	3.28	3 (100%)
2	DMS	B	803	-	3,3,3	0.28	0	3,3,3	3.29	3 (100%)
2	DMS	B	801	-	3,3,3	0.35	0	3,3,3	3.25	3 (100%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	803	DMS	C2-S-C1	4.28	120.35	98.42
2	A	802	DMS	C2-S-C1	4.20	119.94	98.42
2	A	804	DMS	C2-S-C1	4.14	119.62	98.42
2	B	801	DMS	C2-S-C1	3.99	118.89	98.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	DMS	O-S-C1	2.87	120.86	106.49
2	A	804	DMS	O-S-C2	2.80	120.49	106.49
2	B	801	DMS	O-S-C2	2.74	120.22	106.49
2	A	802	DMS	O-S-C2	2.73	120.15	106.49
2	B	803	DMS	O-S-C1	2.69	119.94	106.49
2	A	802	DMS	O-S-C1	2.68	119.91	106.49
2	A	804	DMS	O-S-C1	2.68	119.89	106.49
2	B	803	DMS	O-S-C2	2.64	119.70	106.49

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	804	DMS	2	0
3	A	902	DTZ	1	0
2	B	801	DMS	6	0

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	302/306 (98%)	0.12	8 (2%) 57 64	16, 30, 45, 55	0
1	B	304/306 (99%)	0.02	7 (2%) 61 68	16, 26, 47, 58	0
All	All	606/612 (99%)	0.07	15 (2%) 58 65	16, 28, 46, 58	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	ALA	4.4
1	B	2	GLY	3.3
1	A	215	GLY	3.2
1	A	138	GLY	3.2
1	A	139	SER	3.0
1	B	50	LEU	2.9
1	A	87	LEU	2.4
1	A	154	TYR	2.4
1	B	304	THR	2.4
1	B	3	PHE	2.3
1	A	301	SER	2.3
1	A	74	GLN	2.2
1	B	47	GLU	2.1
1	A	140	PHE	2.0
1	B	296	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	DMS	A	804	4/4	0.87	0.16	63,64,64,64	0
2	DMS	B	803	4/4	0.92	0.14	56,57,58,58	0
3	DTZ	A	902	3/3	0.94	0.23	53,53,59,61	0
2	DMS	B	801	4/4	0.95	0.12	44,47,47,47	0
3	DTZ	B	901	3/3	0.95	0.14	50,50,56,57	0
2	DMS	A	802	4/4	0.97	0.10	36,38,38,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.