



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2024 – 03:47 AM EDT

PDB ID : 1M72  
Title : Crystal Structure of Caspase-1 from Spodoptera frugiperda  
Authors : Forsyth, C.M.; Lemongello, D.; Friesen, P.D.; Fisher, A.J.  
Deposited on : 2002-07-18  
Resolution : 2.30 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.20.1  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.003 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

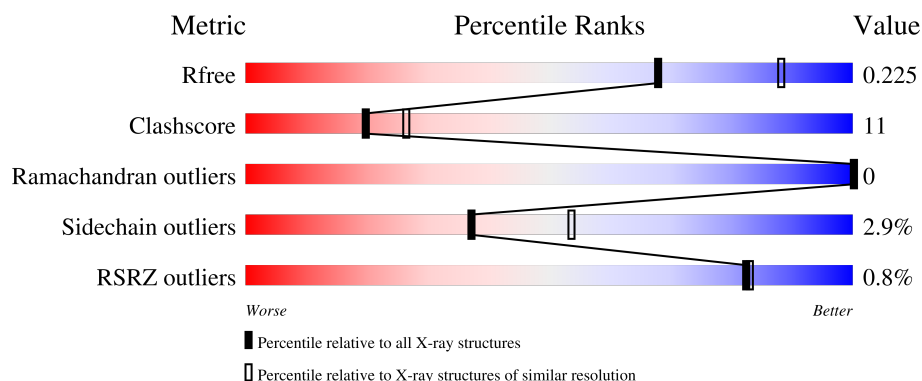
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	164625	5963 (2.30-2.30)
Clashscore	180529	6698 (2.30-2.30)
Ramachandran outliers	177936	6640 (2.30-2.30)
Sidechain outliers	177891	6640 (2.30-2.30)
RSRZ outliers	164620	5963 (2.30-2.30)

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  $\geq 3$ , 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	272	<div> <div></div> <div>68%</div> <div>20%</div> <div>9%</div> </div>
1	B	272	<div> <div></div> <div>73%</div> <div>18%</div> <div>9%</div> </div>
1	C	272	<div> <div></div> <div>75%</div> <div>17%</div> <div>7%</div> </div>
2	D	6	<div> <div></div> <div>67%</div> <div>33%</div> </div>
2	E	6	<div> <div></div> <div>83%</div> <div>17%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	6	 <div>67% 33%</div>

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
3	EDO	B	503	-	-	X	-
3	EDO	C	502	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6433 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 Caspase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1974	1260	341	358	15			
1	B	248	Total	C	N	O	S	0	0	0
			1985	1267	343	360	15			
1	C	252	Total	C	N	O	S	0	0	0
			2016	1286	350	365	15			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	300	LEU	-	cloning artifact	UNP P89116
B	300	LEU	-	cloning artifact	UNP P89116
C	300	LEU	-	cloning artifact	UNP P89116

- Molecule 2 is a protein called Ace-Asp-Glu-Val-Asp-chloromethylketone.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	6	Total	C	N	O	0	0	1
			36	21	4	11			
2	E	6	Total	C	N	O	0	0	1
			36	21	4	11			
2	F	6	Total	C	N	O	0	0	1
			36	21	4	11			

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	102	Total O 102 102	0	0
4	D	1	Total O 1 1	0	0
4	B	92	Total O 92 92	0	0
4	E	4	Total O 4 4	0	0

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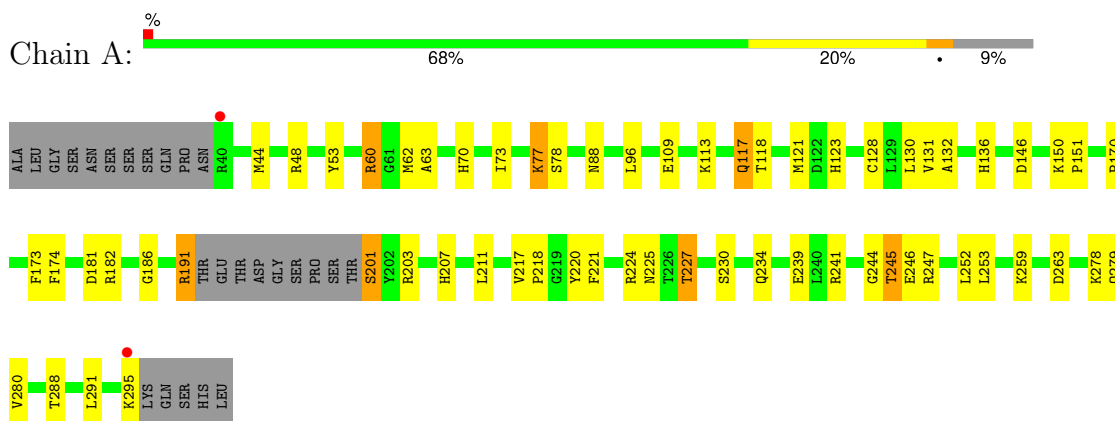
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	116	Total 116	O 116	0	0
4	F	3	Total 3	O 3	0	0

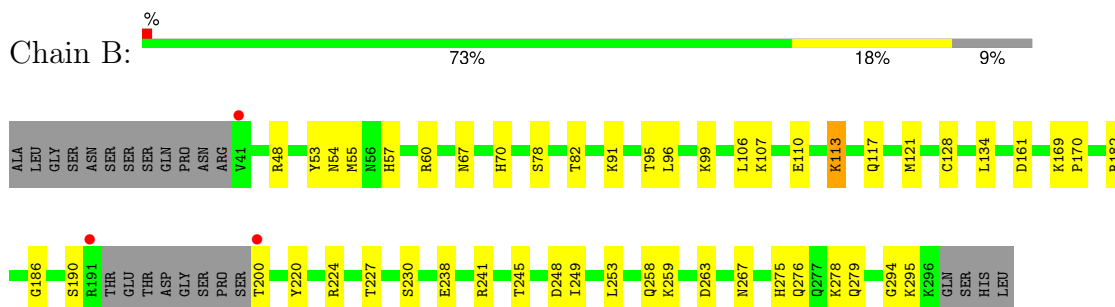
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

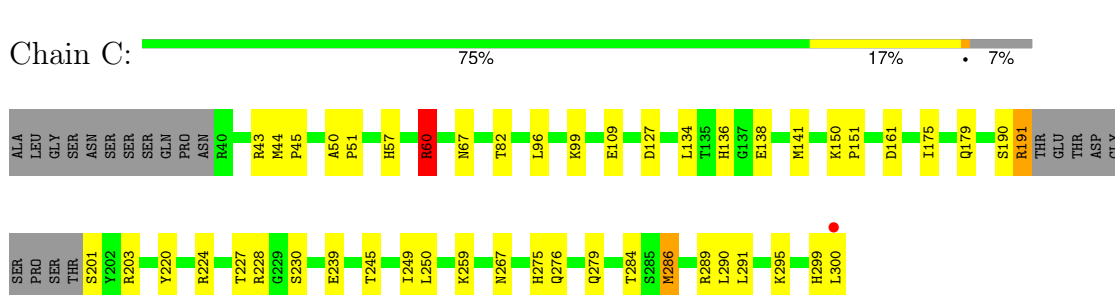
- Molecule 1: Caspase-1



- Molecule 1: Caspase-1



- Molecule 1: Caspase-1




- Molecule 2: Ace-Asp-Glu-Val-Asp-chloromethylketone





- Molecule 2: Ace-Asp-Glu-Val-Asp-chloromethylketone

Chain E:  83% 17%



- Molecule 2: Ace-Asp-Glu-Val-Asp-chloromethylketone

Chain F:  67% 33%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	151.73Å 151.73Å 79.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.30 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.6 (30.00-2.30) 97.6 (30.00-2.30)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.18 (at 2.31Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.183 , 0.232 0.178 , 0.225	Depositor DCC
$R_{free}$ test set	2190 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.4	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6433	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 0QE, ACE, EDO

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	0.74	0/2022	0.86	1/2735 (0.0%)
1	B	0.78	0/2033	0.83	0/2749
1	C	0.80	0/2065	0.89	4/2792 (0.1%)
2	D	1.77	1/32 (3.1%)	0.91	0/43
2	E	1.74	1/32 (3.1%)	0.85	0/43
2	F	1.91	1/32 (3.1%)	0.90	0/43
All	All	0.80	3/6216 (0.0%)	0.86	5/8405 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	500	ACE	C-N	9.98	1.57	1.34
2	E	500	ACE	C-N	9.38	1.55	1.34
2	D	500	ACE	C-N	8.96	1.54	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	60	ARG	NE-CZ-NH2	-8.18	116.21	120.30
1	C	60	ARG	NE-CZ-NH1	6.33	123.46	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	203	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	C	203	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	C	289	ARG	NE-CZ-NH1	5.23	122.91	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	220	TYR	Sidechain
1	B	220	TYR	Sidechain

## 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	1974	0	1938	52	0
1	B	1985	0	1956	38	0
1	C	2016	0	1982	50	0
2	D	36	0	26	1	0
2	E	36	0	26	0	0
2	F	36	0	26	1	0
3	A	4	0	6	2	0
3	B	12	0	18	7	0
3	C	16	0	24	12	0
4	A	102	0	0	3	0
4	B	92	0	0	2	0
4	C	116	0	0	3	0
4	D	1	0	0	0	0
4	E	4	0	0	0	0
4	F	3	0	0	0	0
All	All	6433	0	6002	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 11.

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:62:MET:HE3	1:A:118:THR:HG23	1.31	1.10
1:A:62:MET:CE	1:A:118:THR:HG23	2.02	0.89
1:C:290:LEU:HD23	3:C:502:EDO:H22	1.57	0.86
1:A:280:VAL:H	3:A:506:EDO:H21	1.39	0.83
1:C:275:HIS:CD2	1:C:276:GLN:HE21	1.96	0.83
1:A:77:LYS:HE3	1:A:224:ARG:HH22	1.44	0.82
1:A:191:ARG:HH11	1:A:191:ARG:CB	1.95	0.79
1:C:249:ILE:HG22	3:C:502:EDO:H12	1.65	0.76
1:C:179:GLN:HA	3:C:501:EDO:H11	1.66	0.76
1:C:267:ASN:HD22	1:C:275:HIS:HE1	1.35	0.75
1:C:228:ARG:HD2	4:C:1279:HOH:O	1.87	0.74
1:C:275:HIS:HD2	1:C:276:GLN:HE21	1.34	0.73
1:A:191:ARG:HH11	1:A:191:ARG:HB3	1.54	0.73
1:A:295:LYS:HG3	4:A:1220:HOH:O	1.90	0.72
1:C:43:ARG:HH12	1:C:57:HIS:HE1	1.38	0.72
1:C:60:ARG:HD2	1:C:96:LEU:O	1.90	0.71
1:A:62:MET:HE2	1:A:123:HIS:HE1	1.57	0.70
1:A:77:LYS:HD2	1:A:78:SER:O	1.94	0.67
1:A:60:ARG:HD2	1:A:96:LEU:O	1.94	0.67
1:B:245:THR:HG21	1:B:295:LYS:HG2	1.78	0.66
1:B:117:GLN:O	1:B:121:MET:HG3	1.95	0.66
1:B:113:LYS:O	1:B:117:GLN:HG3	1.95	0.65
1:A:230:SER:O	1:A:234:GLN:HG2	1.96	0.65
1:C:224:ARG:HA	1:C:230:SER:HA	1.80	0.64
1:A:62:MET:CE	1:A:123:HIS:HE1	2.10	0.63
1:B:279:GLN:HA	3:B:507:EDO:H12	1.81	0.63
1:B:275:HIS:CD2	1:B:276:GLN:HG3	2.33	0.62
1:C:267:ASN:HD22	1:C:275:HIS:CE1	2.18	0.62
1:C:245:THR:HB	4:C:1204:HOH:O	1.98	0.62
1:B:54:ASN:HB3	3:B:503:EDO:H12	1.81	0.61
1:B:60:ARG:HD2	1:B:96:LEU:O	2.00	0.60
1:A:224:ARG:HA	1:A:230:SER:HA	1.82	0.60
1:A:88:ASN:OD1	1:A:241:ARG:NH2	2.34	0.60
1:B:267:ASN:HD22	1:B:275:HIS:HE1	1.50	0.60
1:C:295:LYS:NZ	1:C:299:HIS:HD2	2.00	0.60
1:C:99:LYS:HD3	4:C:1253:HOH:O	2.01	0.59
1:B:70:HIS:HD2	1:B:78:SER:OG	1.85	0.59
1:A:245:THR:HG22	1:A:246:GLU:HG3	1.84	0.59
1:A:44:MET:CE	1:A:201:SER:HB3	2.33	0.58
1:C:275:HIS:CD2	1:C:276:GLN:NE2	2.71	0.58
1:B:48:ARG:NH1	4:B:1236:HOH:O	2.37	0.58
1:B:106:LEU:HB3	1:B:110:GLU:HB2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:THR:HG23	1:B:294:GLY:O	2.05	0.57
1:A:279:GLN:HG3	3:A:506:EDO:H22	1.88	0.56
1:C:249:ILE:HG22	3:C:502:EDO:C1	2.35	0.56
1:A:70:HIS:HD2	1:A:78:SER:OG	1.88	0.56
1:A:182:ARG:HH11	1:A:182:ARG:HG2	1.71	0.56
1:A:77:LYS:CE	1:A:224:ARG:HH22	2.17	0.56
1:C:267:ASN:ND2	1:C:275:HIS:HE1	2.01	0.55
1:C:179:GLN:HG2	3:C:501:EDO:H12	1.88	0.55
1:B:91:LYS:O	1:B:95:THR:HG23	2.05	0.55
1:B:267:ASN:ND2	1:B:275:HIS:HE1	2.05	0.54
1:A:217:VAL:HG13	1:A:218:PRO:HD2	1.89	0.54
1:A:48:ARG:NH2	1:B:263:ASP:OD1	2.41	0.54
1:C:179:GLN:HG2	3:C:501:EDO:C1	2.39	0.53
1:A:62:MET:HE3	1:A:118:THR:CG2	2.22	0.53
1:C:44:MET:CE	1:C:201:SER:HB2	2.39	0.53
1:A:136:HIS:HA	2:D:504:ASP:O	2.09	0.52
1:C:191:ARG:NH1	1:C:191:ARG:HG3	2.24	0.52
1:B:224:ARG:HA	1:B:230:SER:HA	1.92	0.51
1:C:191:ARG:CG	1:C:191:ARG:HH11	2.24	0.51
1:A:263:ASP:OD1	1:B:48:ARG:NH2	2.44	0.50
1:A:191:ARG:HH11	1:A:191:ARG:CG	2.25	0.50
1:B:53:TYR:HB3	3:B:503:EDO:H21	1.94	0.49
1:B:67:ASN:ND2	1:B:134:LEU:HD12	2.28	0.49
1:C:295:LYS:NZ	1:C:299:HIS:CD2	2.80	0.49
1:A:109:GLU:HG3	4:A:1267:HOH:O	2.12	0.49
1:B:182:ARG:HG2	1:B:182:ARG:HH11	1.78	0.49
1:C:299:HIS:O	1:C:300:LEU:HB2	2.13	0.49
1:B:99:LYS:HE3	4:B:1240:HOH:O	2.13	0.49
1:C:190:SER:O	1:C:191:ARG:HB2	2.13	0.48
1:A:207:HIS:HE1	4:A:1225:HOH:O	1.97	0.48
1:C:138:GLU:HB2	1:C:141:MET:HG2	1.94	0.48
1:A:239:GLU:OE2	1:A:247:ARG:NH1	2.47	0.48
1:C:239:GLU:OE1	1:C:259:LYS:HD2	2.15	0.47
1:B:53:TYR:HB3	3:B:503:EDO:C2	2.44	0.47
1:B:128:CYS:HB3	1:B:170:PRO:HG2	1.98	0.46
1:B:245:THR:CG2	1:B:295:LYS:HG2	2.45	0.46
1:A:130:LEU:C	1:A:130:LEU:HD23	2.36	0.46
1:B:57:HIS:CE1	1:B:169:LYS:HE2	2.51	0.46
1:B:238:GLU:HA	1:B:241:ARG:NH1	2.30	0.46
1:C:67:ASN:ND2	1:C:134:LEU:HD12	2.30	0.46
1:A:211:LEU:HD21	1:A:253:LEU:CD1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:175:ILE:HD12	1:C:175:ILE:N	2.32	0.45
1:A:77:LYS:HE2	1:A:136:HIS:HE1	1.80	0.45
1:A:288:THR:O	1:B:258:GLN:HB2	2.17	0.45
1:A:253:LEU:HD23	1:A:253:LEU:HA	1.83	0.45
1:C:191:ARG:HG3	1:C:191:ARG:HH11	1.82	0.45
1:A:132:ALA:HA	1:A:174:PHE:O	2.17	0.45
1:C:279:GLN:HA	3:C:504:EDO:H12	1.99	0.44
1:C:220:TYR:CD1	1:C:220:TYR:N	2.85	0.44
1:A:73:ILE:HD11	1:A:146:ASP:C	2.38	0.44
1:B:54:ASN:H	3:B:503:EDO:C1	2.30	0.44
1:A:191:ARG:CG	1:A:191:ARG:NH1	2.81	0.44
1:B:54:ASN:N	3:B:503:EDO:H21	2.33	0.44
1:B:186:GLY:HA3	1:B:278:LYS:HG3	2.00	0.44
1:C:150:LYS:HA	1:C:151:PRO:HD3	1.87	0.44
1:A:131:VAL:O	1:A:173:PHE:HA	2.18	0.43
1:C:43:ARG:HH12	1:C:57:HIS:CE1	2.26	0.43
1:C:161:ASP:OD1	1:C:161:ASP:N	2.49	0.43
1:C:291:LEU:H	3:C:502:EDO:H11	1.83	0.43
1:C:57:HIS:CD2	1:C:127:ASP:HA	2.53	0.43
1:C:249:ILE:CG2	3:C:502:EDO:H12	2.43	0.43
1:A:225:ASN:OD1	1:A:227:THR:HG23	2.17	0.43
1:B:55:MET:HB3	1:B:60:ARG:HH21	1.83	0.43
1:C:67:ASN:ND2	1:C:82:THR:OG1	2.49	0.43
1:B:107:LYS:HB2	1:B:110:GLU:HG3	2.00	0.43
1:C:136:HIS:HA	2:F:504:ASP:O	2.19	0.43
1:C:279:GLN:HA	3:C:504:EDO:C2	2.49	0.43
1:C:295:LYS:HZ1	1:C:299:HIS:HD2	1.66	0.43
1:A:62:MET:CE	1:A:123:HIS:CE1	2.97	0.43
1:A:186:GLY:HA3	1:A:278:LYS:HG3	2.01	0.43
1:C:284:THR:HG22	1:C:286:MET:CE	2.49	0.42
1:A:113:LYS:HE2	1:A:113:LYS:HB3	1.79	0.42
1:C:109:GLU:H	1:C:109:GLU:CD	2.22	0.42
1:B:161:ASP:N	1:B:161:ASP:OD1	2.52	0.42
1:A:128:CYS:HB2	1:A:170:PRO:O	2.19	0.42
1:B:54:ASN:H	3:B:503:EDO:H12	1.84	0.42
1:B:67:ASN:ND2	1:B:82:THR:OG1	2.51	0.42
1:C:50:ALA:HA	1:C:51:PRO:HD3	1.83	0.42
1:C:191:ARG:NH1	1:C:191:ARG:CG	2.80	0.42
1:A:117:GLN:O	1:A:121:MET:HG3	2.20	0.41
1:A:181:ASP:O	1:A:221:PHE:CE1	2.73	0.41
1:A:182:ARG:HG2	1:A:182:ARG:NH1	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:GLU:OE1	1:A:259:LYS:HE3	2.19	0.41
1:B:253:LEU:HD23	1:B:253:LEU:HA	1.84	0.41
1:A:44:MET:HE1	1:A:201:SER:HB3	1.99	0.41
1:B:249:ILE:HD12	1:B:249:ILE:HA	1.87	0.41
1:A:53:TYR:CD1	1:A:170:PRO:HD3	2.56	0.41
1:B:55:MET:HB3	1:B:60:ARG:NH2	2.35	0.41
1:A:63:ALA:HA	1:A:130:LEU:O	2.20	0.41
1:C:250:LEU:HG	3:C:502:EDO:H21	2.03	0.41
1:C:67:ASN:HD22	1:C:134:LEU:HD12	1.85	0.40
1:C:275:HIS:CD2	1:C:276:GLN:HG3	2.57	0.40
1:A:150:LYS:HA	1:A:151:PRO:HD3	1.90	0.40
1:A:244:GLY:HA2	1:A:252:LEU:HD11	2.02	0.40
1:C:290:LEU:HA	3:C:502:EDO:C2	2.52	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	243/272 (89%)	234 (96%)	9 (4%)	0	100	100
1	B	244/272 (90%)	238 (98%)	6 (2%)	0	100	100
1	C	248/272 (91%)	241 (97%)	7 (3%)	0	100	100
2	D	3/6 (50%)	3 (100%)	0	0	100	100
2	E	3/6 (50%)	3 (100%)	0	0	100	100
2	F	3/6 (50%)	3 (100%)	0	0	100	100
All	All	744/834 (89%)	722 (97%)	22 (3%)	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	215/239 (90%)	207 (96%)	8 (4%)	29	43
1	B	217/239 (91%)	211 (97%)	6 (3%)	38	55
1	C	220/239 (92%)	215 (98%)	5 (2%)	45	63
2	D	4/4 (100%)	4 (100%)	0	100	100
2	E	4/4 (100%)	4 (100%)	0	100	100
2	F	4/4 (100%)	4 (100%)	0	100	100
All	All	664/729 (91%)	645 (97%)	19 (3%)	37	54

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ARG
1	A	77	LYS
1	A	117	GLN
1	A	191	ARG
1	A	201	SER
1	A	227	THR
1	A	245	THR
1	A	291	LEU
1	B	113	LYS
1	B	190	SER
1	B	200	THR
1	B	227	THR
1	B	248	ASP
1	B	259	LYS
1	C	45	PRO
1	C	60	ARG
1	C	191	ARG
1	C	227	THR
1	C	286	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	67	ASN
1	A	70	HIS
1	A	123	HIS
1	A	153	ASN
1	A	267	ASN
1	A	276	GLN
1	B	57	HIS
1	B	67	ASN
1	B	70	HIS
1	B	83	ASN
1	B	153	ASN
1	B	258	GLN
1	B	267	ASN
1	B	276	GLN
1	C	57	HIS
1	C	67	ASN
1	C	83	ASN
1	C	153	ASN
1	C	267	ASN
1	C	275	HIS
1	C	299	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

8 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	$\# Z  > 2$
3	EDO	B	505	-	3,3,3	0.68	0	2,2,2	0.26	0
3	EDO	C	504	-	3,3,3	0.42	0	2,2,2	0.59	0
3	EDO	B	503	-	3,3,3	0.20	0	2,2,2	0.50	0
3	EDO	C	501	-	3,3,3	0.63	0	2,2,2	0.68	0
3	EDO	A	506	-	3,3,3	0.41	0	2,2,2	0.39	0
3	EDO	C	508	-	3,3,3	0.63	0	2,2,2	0.33	0
3	EDO	B	507	-	3,3,3	0.37	0	2,2,2	0.60	0
3	EDO	C	502	-	3,3,3	0.23	0	2,2,2	0.46	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
3	EDO	B	505	-	-	1/1/1/1	-
3	EDO	C	504	-	-	0/1/1/1	-
3	EDO	B	503	-	-	1/1/1/1	-
3	EDO	C	501	-	-	0/1/1/1	-
3	EDO	A	506	-	-	0/1/1/1	-
3	EDO	C	508	-	-	0/1/1/1	-
3	EDO	B	507	-	-	0/1/1/1	-
3	EDO	C	502	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	505	EDO	O1-C1-C2-O2
3	B	503	EDO	O1-C1-C2-O2
3	C	502	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	504	EDO	2	0
3	B	503	EDO	6	0
3	C	501	EDO	3	0
3	A	506	EDO	2	0
3	B	507	EDO	1	0
3	C	502	EDO	7	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	247/272 (90%)	-0.41	2 (0%) 82 83	18, 29, 46, 62	0
1	B	248/272 (91%)	-0.49	3 (1%) 76 76	16, 28, 45, 67	0
1	C	252/272 (92%)	-0.64	1 (0%) 89 89	13, 24, 40, 60	0
2	D	4/6 (66%)	-0.13	0 100 100	34, 37, 39, 41	0
2	E	4/6 (66%)	-0.31	0 100 100	29, 36, 40, 40	0
2	F	4/6 (66%)	-0.21	0 100 100	30, 32, 37, 40	0
All	All	759/834 (91%)	-0.51	6 (0%) 82 83	13, 27, 45, 67	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	300	LEU	3.7
1	A	40	ARG	2.9
1	B	41	VAL	2.7
1	B	191	ARG	2.7
1	A	295	LYS	2.7
1	B	200	THR	2.3

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	EDO	B	505	4/4	0.92	0.11	33,41,42,42	0
3	EDO	B	507	4/4	0.93	0.15	35,35,36,37	0
3	EDO	C	501	4/4	0.93	0.14	24,27,29,32	0
3	EDO	C	508	4/4	0.93	0.11	45,48,48,48	0
3	EDO	A	506	4/4	0.94	0.13	31,34,35,36	0
3	EDO	C	504	4/4	0.95	0.12	37,40,40,42	0
3	EDO	B	503	4/4	0.96	0.15	24,27,30,30	0
3	EDO	C	502	4/4	0.97	0.09	28,29,30,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.