



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:58 PM GMT

PDB ID : 1I4O  
Title : CRYSTAL STRUCTURE OF THE XIAP/CASPASE-7 COMPLEX  
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Deposited on : 2001-02-22  
Resolution : 2.40 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

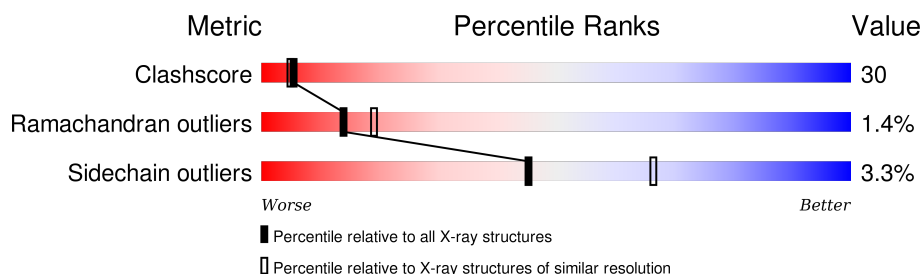
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	280	
1	B	280	
2	C	141	
2	D	141	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4124 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-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	0	0
			1854	1181	313	345	15			
1	B	232	Total	C	N	O	S	0	0	0
			1823	1161	310	337	15			

- Molecule 2 is a protein called BACULOVIRAL IAP REPEAT-CONTAINING PROTEIN 4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	17	Total	C	N	O	0	0	0
			124	78	20	26			
2	D	17	Total	C	N	O	0	0	0
			124	78	20	26			

- Molecule 3 is water.

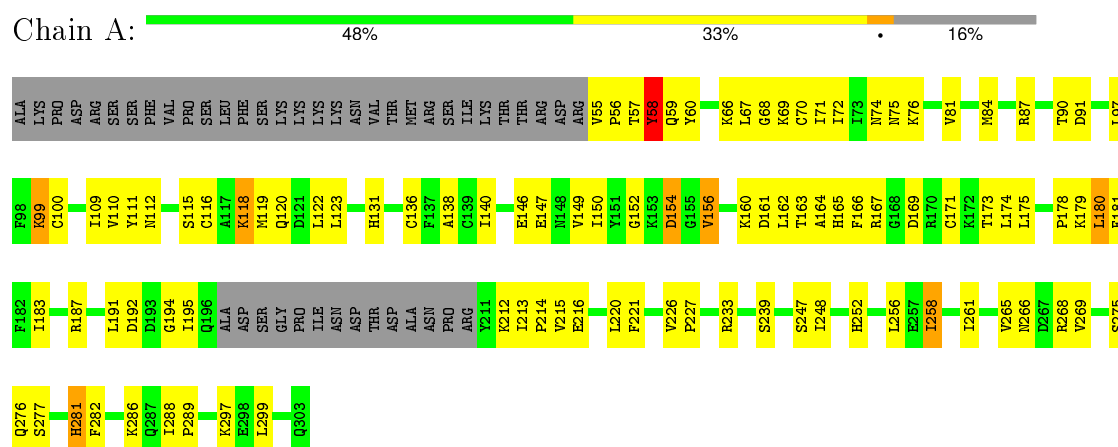
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	87	Total	O	0	0
			87	87		
3	B	98	Total	O	0	0
			98	98		
3	C	6	Total	O	0	0
			6	6		
3	D	8	Total	O	0	0
			8	8		

### 3 Residue-property plots

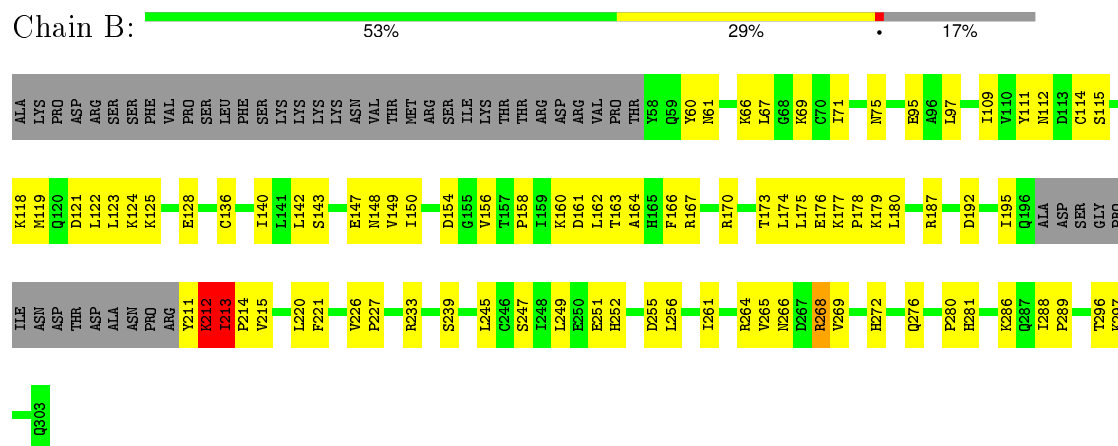
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

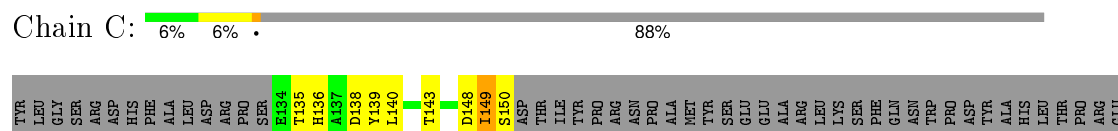
#### • Molecule 1: CASPASE-7



#### • Molecule 1: CASPASE-7



#### • Molecule 2: BACULOVIRAL IAP REPEAT-CONTAINING PROTEIN 4



LEU ALA SER ALA GLY LEU TYR TYR THR GLY ILE GLY ASP GLN VAL GLN CYS CYS PHE CYS CYS GLY GLY LYS LEU LYS ASN TRP TRP GLU PRO CYS ASP ARG ALA TRP TRP SER LEU HIS ARG ARG ARG HIS PHE PRO ASN CYS PHE VAL LEU GLY ARG ASN ASN ILE ARG SER GLU SER ASP ALA

VAL  
SER  
SER  
ASP  
ARG  
ASN  
PHE  
PRO  
ASN  
SER  
THR  
ASN  
LEU  
PRO  
ARG  
ASN  
PRO

- Molecule 2: BACULOVIRAL IAP REPEAT-CONTAINING PROTEIN 4

Chain D:  6% 6% • 88%

Tyr	Leu	Gly	Ser	Arg	Asp	His	Phe	Leu	Leu	Asp	Pro	Ser	Ala	His	Met	Ser	Glu	Glu	Ala	Arg	Leu	Lys	Ser	Phe	Gln	Asn	Trp	Pro	Asp	Tyr	Ala	His	Leu	Thr	Pro	Asp
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

[illegible]

ALA  
VAL  
SER  
SER  
ASP  
ARG  
ASN  
PHE  
PRO  
ASN  
SER  
THR  
ASN  
LEU  
PRO  
ARG  
ASN  
PRO

## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.50Å 88.50Å 185.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.40)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	?	Depositor
R, $R_{free}$	0.224 , 0.260	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4124	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.44	0/1895	0.73	2/2556 (0.1%)
1	B	0.44	0/1862	0.72	2/2510 (0.1%)
2	C	0.77	0/125	0.68	0/171
2	D	0.69	0/125	0.64	0/171
All	All	0.47	0/4007	0.72	4/5408 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	213	ILE	N-CA-C	-6.64	93.08	111.00
1	A	152	GLY	N-CA-C	-5.87	98.42	113.10
1	A	58	TYR	N-CA-C	5.08	124.71	111.00
1	B	212	LYS	N-CA-C	-5.05	97.36	111.00

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	1854	0	1793	125	0
1	B	1823	0	1763	106	0
2	C	124	0	113	15	0
2	D	124	0	113	16	0
3	A	87	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	98	0	0	19	0
3	C	6	0	0	2	0
3	D	8	0	0	1	0
All	All	4124	0	3782	230	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 30.

All (230) 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:194:GLY:HA2	1:B:211:TYR:CB	1.87	1.04
1:B:75:ASN:HD21	1:B:119:MET:HE1	1.25	0.99
1:A:276:GLN:O	2:C:149:ILE:HG23	1.63	0.99
1:A:286:LYS:HG3	1:B:212:LYS:H	1.30	0.95
2:C:149:ILE:HG12	2:C:149:ILE:O	1.71	0.90
1:B:75:ASN:HD21	1:B:119:MET:CE	1.85	0.89
2:D:136:HIS:O	2:D:140:LEU:HD23	1.75	0.86
1:B:252:HIS:HB3	1:B:256:LEU:HG	1.61	0.83
3:B:341:HOH:O	2:D:137:ALA:HB2	1.79	0.83
1:B:75:ASN:HD22	1:B:143:SER:HB2	1.43	0.82
1:A:97:LEU:HD13	1:A:140:ILE:HG21	1.61	0.82
1:A:258:ILE:HD11	1:A:299:LEU:HG	1.62	0.79
1:A:269:VAL:HG11	1:A:289:PRO:CG	2.13	0.79
2:D:149:ILE:O	2:D:149:ILE:HG12	1.84	0.78
1:A:81:VAL:HG23	3:A:360:HOH:O	1.84	0.77
2:D:135:THR:HG23	2:D:138:ASP:HB2	1.66	0.77
1:B:111:TYR:CZ	1:B:122:LEU:HD11	2.21	0.76
1:A:175:LEU:HD22	1:A:213:ILE:HD11	1.67	0.76
1:A:233:ARG:HA	1:A:239:SER:HA	1.67	0.76
1:A:194:GLY:HA2	1:B:211:TYR:CA	2.17	0.74
1:B:233:ARG:HA	1:B:239:SER:HA	1.68	0.74
1:A:258:ILE:CG1	1:A:299:LEU:HB3	2.19	0.73
1:B:276:GLN:O	2:D:149:ILE:HG22	1.89	0.73
1:B:211:TYR:HA	1:B:213:ILE:O	1.89	0.73
1:B:211:TYR:C	1:B:213:ILE:H	1.80	0.72
1:B:269:VAL:HG21	1:B:289:PRO:CG	2.19	0.72
1:A:194:GLY:CA	1:B:211:TYR:CB	2.65	0.72
1:B:211:TYR:C	1:B:213:ILE:N	2.39	0.72
1:A:277:SER:HA	2:C:149:ILE:HG21	1.73	0.70
1:A:149:VAL:HG11	1:A:156:VAL:HG13	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:GLU:HG2	3:B:400:HOH:O	1.92	0.69
1:A:269:VAL:HG11	1:A:289:PRO:HG2	1.74	0.68
2:C:139:TYR:HA	3:C:106:HOH:O	1.92	0.68
1:A:74:ASN:HB3	1:A:112:ASN:ND2	2.08	0.68
1:A:55:VAL:N	1:A:56:PRO:O	2.26	0.68
1:A:123:LEU:HD12	1:A:162:LEU:HD22	1.76	0.68
1:A:149:VAL:CG1	1:A:156:VAL:HG13	2.24	0.67
1:A:266:ASN:OD1	1:A:289:PRO:HB2	1.93	0.67
1:A:163:THR:HG21	1:A:221:PHE:HE2	1.59	0.67
1:A:194:GLY:C	1:B:211:TYR:N	2.48	0.67
1:A:55:VAL:HA	1:B:252:HIS:NE2	2.10	0.67
1:B:164:ALA:O	1:B:167:ARG:HG3	1.95	0.66
1:B:213:ILE:O	1:B:213:ILE:HG23	1.94	0.66
1:A:216:GLU:HG3	3:A:316:HOH:O	1.95	0.66
1:A:163:THR:HG21	1:A:221:PHE:CE2	2.31	0.66
1:B:160:LYS:HD2	3:B:394:HOH:O	1.95	0.66
1:A:149:VAL:HG11	1:A:156:VAL:CG1	2.27	0.65
1:B:226:VAL:HG13	1:B:227:PRO:HD2	1.78	0.64
2:D:136:HIS:CE1	2:D:140:LEU:HD21	2.32	0.64
1:A:115:SER:HB2	1:A:154:ASP:OD2	1.97	0.64
1:A:258:ILE:HG13	1:A:299:LEU:HD23	1.80	0.64
1:B:255:ASP:HB2	3:B:367:HOH:O	1.98	0.63
1:A:136:CYS:HB3	1:A:178:PRO:HG2	1.81	0.63
1:A:58:TYR:O	1:A:297:LYS:HB3	1.98	0.63
1:B:211:TYR:HA	1:B:213:ILE:C	2.20	0.62
1:B:97:LEU:HD13	1:B:140:ILE:HG21	1.81	0.62
1:B:147:GLU:HG2	3:B:324:HOH:O	1.97	0.62
1:A:258:ILE:HG12	1:A:299:LEU:HB3	1.79	0.62
1:A:120:GLN:HE22	1:A:161:ASP:HB3	1.64	0.62
1:B:296:THR:O	1:B:297:LYS:HG3	1.99	0.62
1:B:75:ASN:ND2	1:B:119:MET:CE	2.62	0.62
2:C:136:HIS:HB3	3:C:74:HOH:O	2.00	0.61
1:A:261:ILE:O	1:A:265:VAL:HG23	2.00	0.61
1:A:136:CYS:CB	1:A:178:PRO:HG2	2.31	0.61
3:B:332:HOH:O	2:D:149:ILE:HG21	2.01	0.61
1:A:286:LYS:HG3	1:B:211:TYR:CB	2.31	0.61
1:A:165:HIS:HE1	3:A:313:HOH:O	1.84	0.61
1:A:247:SER:HB3	1:A:268:ARG:HH12	1.65	0.60
1:B:136:CYS:HB3	1:B:178:PRO:HG2	1.82	0.60
2:D:135:THR:HG23	2:D:138:ASP:CB	2.31	0.60
1:B:211:TYR:CA	1:B:213:ILE:O	2.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:PRO:HA	3:B:401:HOH:O	2.01	0.59
1:A:118:LYS:HD3	3:A:386:HOH:O	2.02	0.59
1:B:266:ASN:OD1	1:B:289:PRO:HB2	2.02	0.59
1:B:150:ILE:O	1:B:150:ILE:HD12	2.02	0.59
1:A:75:ASN:HB3	3:A:375:HOH:O	2.03	0.59
1:B:115:SER:HB2	1:B:154:ASP:OD2	2.03	0.58
1:B:272:HIS:ND1	1:B:272:HIS:O	2.37	0.58
1:A:297:LYS:HE3	3:B:342:HOH:O	2.02	0.58
1:A:277:SER:HA	2:C:149:ILE:CG2	2.34	0.58
1:B:166:PHE:HB3	1:B:179:LYS:HD2	1.85	0.58
1:A:120:GLN:HB3	3:A:313:HOH:O	2.03	0.58
1:B:69:LYS:HE2	1:B:71:ILE:HD11	1.85	0.57
1:B:149:VAL:HB	1:B:156:VAL:HG22	1.87	0.57
1:B:124:LYS:O	1:B:128:GLU:HG3	2.04	0.57
1:A:69:LYS:HE2	1:A:71:ILE:HD11	1.88	0.56
2:C:148:ASP:OD1	2:C:148:ASP:O	2.24	0.56
1:A:220:LEU:HD13	1:A:221:PHE:N	2.22	0.55
1:A:183:ILE:N	1:A:183:ILE:HD12	2.22	0.55
1:A:76:LYS:HD2	1:A:91:ASP:OD2	2.07	0.55
1:A:147:GLU:HB2	1:A:187:ARG:O	2.08	0.54
1:B:212:LYS:HD2	1:B:212:LYS:N	2.23	0.54
1:A:120:GLN:OE1	1:A:162:LEU:HD23	2.08	0.54
1:B:220:LEU:C	1:B:220:LEU:HD23	2.27	0.54
1:A:118:LYS:HA	1:A:118:LYS:HE3	1.89	0.54
1:A:99:LYS:HD2	1:A:100:CYS:N	2.22	0.54
1:A:248:ILE:HG22	1:A:261:ILE:HG23	1.89	0.54
1:B:147:GLU:O	1:B:148:ASN:HB2	2.08	0.53
1:A:195:ILE:HG22	1:B:211:TYR:N	2.23	0.53
1:A:247:SER:CB	1:A:268:ARG:NH1	2.72	0.53
1:A:226:VAL:HG13	1:A:227:PRO:HD2	1.91	0.53
1:A:276:GLN:HB3	2:C:150:SER:HB3	1.90	0.52
1:B:160:LYS:HG2	3:B:329:HOH:O	2.10	0.52
1:B:245:LEU:O	1:B:249:LEU:HB2	2.09	0.52
1:A:171:CYS:SG	1:A:174:LEU:HD12	2.50	0.52
1:B:211:TYR:O	1:B:212:LYS:HB2	2.10	0.51
2:D:140:LEU:HD22	2:D:140:LEU:N	2.26	0.51
1:A:258:ILE:HG13	1:A:299:LEU:HB3	1.91	0.51
1:B:269:VAL:HG21	1:B:289:PRO:HG3	1.93	0.51
1:B:60:TYR:CD1	1:B:178:PRO:HD3	2.46	0.51
1:A:150:ILE:HD12	1:A:150:ILE:O	2.10	0.51
1:A:194:GLY:CA	1:B:211:TYR:N	2.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LEU:HD21	1:B:195:ILE:HG21	1.93	0.51
1:A:58:TYR:CG	1:A:59:GLN:N	2.78	0.51
1:A:87:ARG:HG2	1:A:233:ARG:HH12	1.75	0.51
1:A:220:LEU:HD13	1:A:220:LEU:C	2.31	0.51
1:A:252:HIS:HB3	1:A:256:LEU:HG	1.92	0.51
1:A:191:LEU:HD11	1:A:282:PHE:HD1	1.76	0.50
2:D:139:TYR:HD2	2:D:140:LEU:HD22	1.76	0.50
1:B:142:LEU:HB3	3:B:344:HOH:O	2.11	0.50
2:D:143:THR:OG1	2:D:145:GLN:HG3	2.10	0.50
1:A:84:MET:SD	2:C:143:THR:HA	2.52	0.50
1:B:136:CYS:CB	1:B:178:PRO:HG2	2.40	0.50
1:B:112:ASN:HB3	3:B:314:HOH:O	2.12	0.50
1:A:146:GLU:O	1:A:147:GLU:C	2.50	0.49
1:A:99:LYS:C	1:A:99:LYS:HD2	2.32	0.49
1:A:164:ALA:O	1:A:167:ARG:HG3	2.12	0.49
2:D:139:TYR:CD2	2:D:140:LEU:HD22	2.46	0.49
1:A:194:GLY:N	1:B:211:TYR:CB	2.76	0.49
1:A:258:ILE:HD11	1:A:299:LEU:CG	2.37	0.49
1:A:286:LYS:HB3	1:B:214:PRO:HG3	1.94	0.49
1:B:150:ILE:HD11	1:B:162:LEU:HD11	1.95	0.49
1:A:166:PHE:HB3	1:A:179:LYS:HD2	1.93	0.49
1:A:276:GLN:O	1:A:276:GLN:HG2	2.12	0.49
1:A:214:PRO:HG3	1:B:286:LYS:HB3	1.95	0.49
1:A:247:SER:HB3	1:A:268:ARG:NH1	2.28	0.49
1:A:226:VAL:CG1	1:A:227:PRO:HD2	2.44	0.48
1:A:276:GLN:O	2:C:149:ILE:CG2	2.50	0.48
2:C:136:HIS:O	2:C:140:LEU:HD23	2.13	0.48
1:B:281:HIS:CE1	3:B:341:HOH:O	2.66	0.48
1:A:55:VAL:N	1:A:57:THR:HG23	2.29	0.48
1:B:226:VAL:HG13	1:B:227:PRO:CD	2.44	0.48
1:A:215:VAL:HG22	1:B:192:ASP:OD2	2.13	0.47
1:A:195:ILE:N	1:B:211:TYR:O	2.47	0.47
1:A:248:ILE:CG2	1:A:261:ILE:HG23	2.43	0.47
1:A:180:LEU:N	1:A:180:LEU:CD1	2.77	0.47
1:B:272:HIS:HB2	3:B:325:HOH:O	2.13	0.47
2:C:135:THR:HG23	2:C:138:ASP:CG	2.35	0.47
1:B:173:THR:HG23	3:B:393:HOH:O	2.14	0.47
1:B:211:TYR:HA	1:B:214:PRO:CA	2.45	0.47
1:B:69:LYS:HE3	1:B:109:ILE:HD12	1.96	0.47
1:B:158:PRO:CG	1:B:161:ASP:OD2	2.62	0.47
1:A:169:ASP:CB	1:B:195:ILE:HD13	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:THR:HG21	1:B:221:PHE:CE2	2.50	0.47
1:A:212:LYS:HG3	1:B:195:ILE:O	2.15	0.47
1:B:213:ILE:CG2	1:B:213:ILE:O	2.63	0.47
1:A:70:CYS:HA	1:A:138:ALA:O	2.15	0.47
1:A:131:HIS:H	1:A:173:THR:CB	2.28	0.47
1:B:211:TYR:HA	1:B:214:PRO:HA	1.97	0.46
1:B:163:THR:HG21	1:B:221:PHE:CZ	2.51	0.46
1:A:286:LYS:HG3	1:B:212:LYS:N	2.14	0.46
1:A:226:VAL:HG23	1:A:288:ILE:HG23	1.98	0.46
1:A:55:VAL:CA	1:B:252:HIS:NE2	2.78	0.46
1:B:160:LYS:HE3	3:B:370:HOH:O	2.15	0.46
1:A:281:HIS:CD2	3:A:320:HOH:O	2.68	0.46
1:B:211:TYR:N	1:B:215:VAL:HG13	2.31	0.46
1:A:116:CYS:HA	1:A:119:MET:HE2	1.97	0.46
1:A:87:ARG:HH12	1:A:90:THR:CB	2.29	0.46
1:A:149:VAL:HG13	1:A:156:VAL:HG13	1.97	0.46
1:B:269:VAL:CG2	1:B:289:PRO:HD3	2.46	0.45
1:B:288:ILE:HD12	1:B:289:PRO:CD	2.47	0.45
1:B:160:LYS:CE	3:B:370:HOH:O	2.63	0.45
1:A:66:LYS:HE2	3:A:368:HOH:O	2.15	0.45
2:D:145:GLN:NE2	3:D:263:HOH:O	2.47	0.45
1:A:258:ILE:HD13	1:A:258:ILE:N	2.30	0.45
1:A:226:VAL:CG1	1:A:227:PRO:CD	2.95	0.45
1:A:55:VAL:N	1:A:56:PRO:C	2.70	0.44
1:B:288:ILE:HD12	1:B:289:PRO:HD2	1.99	0.44
1:A:60:TYR:HA	3:A:380:HOH:O	2.18	0.44
1:B:69:LYS:HE3	1:B:109:ILE:CD1	2.48	0.44
2:D:149:ILE:O	2:D:149:ILE:CG1	2.58	0.44
1:B:150:ILE:C	1:B:150:ILE:HD12	2.37	0.44
1:B:69:LYS:CE	1:B:109:ILE:HD12	2.47	0.44
1:A:192:ASP:HB3	1:A:286:LYS:O	2.17	0.44
1:B:247:SER:O	1:B:251:GLU:HG3	2.18	0.44
1:A:111:TYR:CG	1:A:122:LEU:HD21	2.53	0.44
1:A:72:ILE:HB	1:A:110:VAL:HG22	2.00	0.44
1:A:149:VAL:CG1	1:A:150:ILE:N	2.81	0.43
1:A:149:VAL:HG12	1:A:150:ILE:N	2.34	0.43
1:A:226:VAL:HG13	1:A:227:PRO:CD	2.48	0.43
2:C:136:HIS:CE1	2:C:140:LEU:HD21	2.53	0.43
1:A:166:PHE:O	1:A:167:ARG:C	2.56	0.43
1:A:123:LEU:CD1	1:A:162:LEU:HB3	2.49	0.43
2:D:135:THR:HG23	2:D:138:ASP:CG	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:LYS:O	1:B:286:LYS:HE2	2.19	0.43
1:A:163:THR:HG22	1:A:181:PHE:CE2	2.54	0.43
1:A:118:LYS:O	1:A:122:LEU:HB2	2.19	0.43
2:D:140:LEU:N	2:D:140:LEU:CD2	2.82	0.42
1:B:261:ILE:O	1:B:265:VAL:HG23	2.18	0.42
1:A:195:ILE:O	1:B:211:TYR:O	2.36	0.42
1:B:175:LEU:O	1:B:176:GLU:HB2	2.19	0.42
1:A:136:CYS:HB2	1:A:178:PRO:HG2	1.99	0.42
2:C:148:ASP:O	2:C:149:ILE:C	2.57	0.42
1:B:226:VAL:CG1	1:B:227:PRO:HD2	2.49	0.42
1:B:170:ARG:HH11	1:B:170:ARG:HG3	1.84	0.42
1:A:123:LEU:HD12	1:A:162:LEU:CD2	2.48	0.42
1:A:178:PRO:HB2	1:A:180:LEU:HD11	2.02	0.42
1:B:288:ILE:HA	1:B:289:PRO:HD3	1.94	0.42
1:B:114:CYS:HB3	1:B:118:LYS:HB3	2.02	0.42
1:A:57:THR:OG1	1:B:264:ARG:HG2	2.20	0.41
1:B:187:ARG:NH2	1:B:227:PRO:HG3	2.36	0.41
1:A:216:GLU:HB3	1:B:288:ILE:HG21	2.03	0.41
1:A:74:ASN:HB3	1:A:112:ASN:HD22	1.83	0.41
1:A:265:VAL:O	1:A:269:VAL:HG12	2.21	0.41
1:A:58:TYR:N	1:A:58:TYR:CD1	2.86	0.41
1:B:147:GLU:HG3	1:B:148:ASN:ND2	2.35	0.41
1:B:66:LYS:HB3	3:B:319:HOH:O	2.21	0.41
1:A:286:LYS:HB2	1:B:211:TYR:CB	2.50	0.41
1:A:226:VAL:CG2	1:A:288:ILE:HG23	2.51	0.41
1:A:68:GLY:HA3	1:A:136:CYS:O	2.21	0.41
1:A:286:LYS:CG	1:B:211:TYR:CB	2.97	0.41
1:B:269:VAL:HG21	1:B:289:PRO:CD	2.50	0.41
1:A:69:LYS:HE3	1:A:109:ILE:HD12	2.03	0.41
1:B:268:ARG:NE	3:B:358:HOH:O	2.53	0.41
1:B:123:LEU:HD12	1:B:162:LEU:HD22	2.03	0.41
1:B:174:LEU:HA	1:B:177:LYS:HD2	2.02	0.41
1:A:288:ILE:HG23	1:A:288:ILE:O	2.21	0.40
2:C:149:ILE:CG1	2:C:149:ILE:O	2.50	0.40
1:B:255:ASP:CB	3:B:367:HOH:O	2.63	0.40
1:B:121:ASP:OD2	1:B:125:LYS:HE3	2.20	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	231/280 (82%)	210 (91%)	17 (7%)	4 (2%)	11	14
1	B	228/280 (81%)	221 (97%)	6 (3%)	1 (0%)	39	56
2	C	15/141 (11%)	14 (93%)	0	1 (7%)	1	0
2	D	15/141 (11%)	14 (93%)	0	1 (7%)	1	0
All	All	489/842 (58%)	459 (94%)	23 (5%)	7 (1%)	14	19

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	ASP
1	A	58	TYR
1	A	275	SER
1	B	213	ILE
1	A	281	HIS
2	C	149	ILE
2	D	149	ILE

### 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	200/248 (81%)	192 (96%)	8 (4%)	38	58
1	B	195/248 (79%)	189 (97%)	6 (3%)	47	69
2	C	13/124 (10%)	13 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	13/124 (10%)	13 (100%)	0	100	100
All	All	421/744 (57%)	407 (97%)	14 (3%)	45	66

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	TYR
1	A	67	LEU
1	A	99	LYS
1	A	118	LYS
1	A	156	VAL
1	A	160	LYS
1	A	180	LEU
1	A	258	ILE
1	B	61	ASN
1	B	67	LEU
1	B	95	GLU
1	B	180	LEU
1	B	212	LYS
1	B	268	ARG

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

Mol	Chain	Res	Type
1	A	74	ASN
1	A	112	ASN
1	B	59	GLN
1	B	61	ASN
1	B	63	ASN
1	B	74	ASN
1	B	75	ASN
1	B	112	ASN
1	B	148	ASN
2	C	145	GLN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.