



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:10 PM GMT

PDB ID : 1EAI  
Title : COMPLEX OF ASCARIS CHYMOTRPSIN/ELASTASE INHIBITOR  
WITH PORCINE ELASTASE  
Authors : Huang, K.; Strynadka, N.C.J.; Bernard, V.D.; Peanasky, R.J.; James, M.N.G.  
Deposited on : 1999-03-25  
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

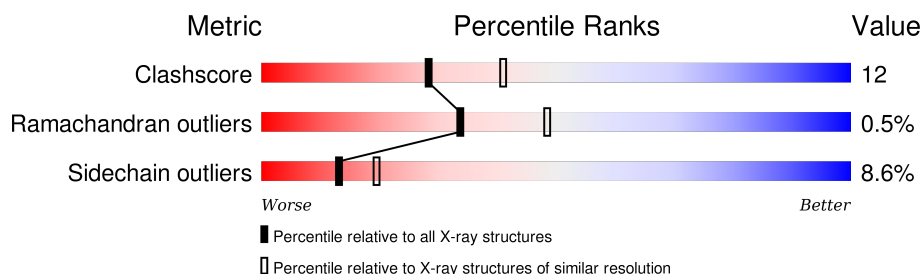
# 1 Overall quality at a glance

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	240	
1	B	240	
2	C	61	
2	D	61	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4686 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 PROTEIN (ELASTASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1822	1135	329	348	10			
1	B	240	Total	C	N	O	S	0	0	0
			1822	1135	329	348	10			

- Molecule 2 is a protein called PROTEIN (CHYMOTRYPSIN/ELASTASE ISOINHIBITOR 1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	61	Total	C	N	O	S	0	0	0
			448	260	84	91	13			
2	D	61	Total	C	N	O	S	0	0	0
			448	260	84	91	13			

- Molecule 3 is water.

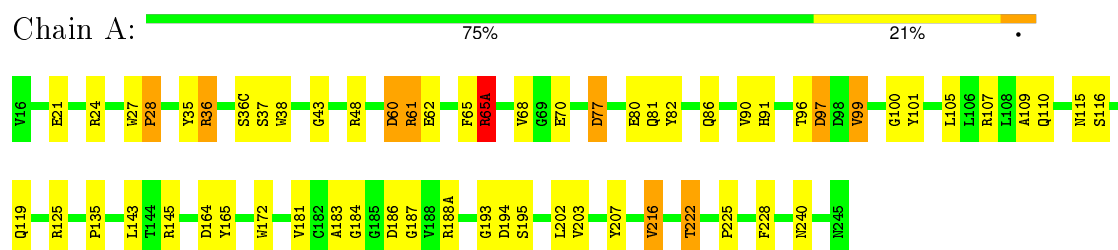
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	58	Total	O	0	0
			58	58		
3	B	61	Total	O	0	0
			61	61		
3	C	16	Total	O	0	0
			16	16		
3	D	11	Total	O	0	0
			11	11		

### 3 Residue-property plots

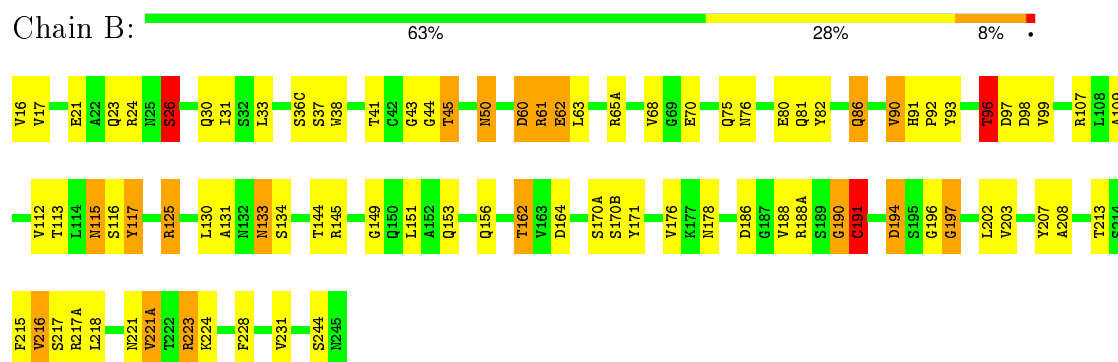
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

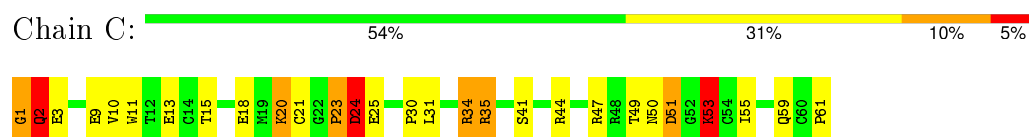
#### • Molecule 1: PROTEIN (ELASTASE)



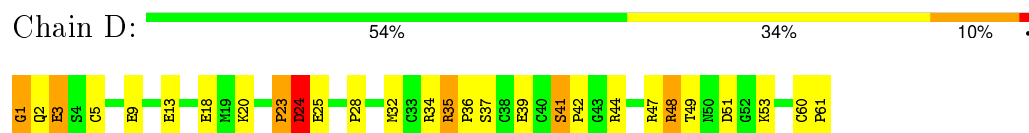
#### • Molecule 1: PROTEIN (ELASTASE)



#### • Molecule 2: PROTEIN (CHYMOTRYPSIN/ELASTASE ISOINHIBITOR 1)



#### • Molecule 2: PROTEIN (CHYMOTRYPSIN/ELASTASE ISOINHIBITOR 1)



## 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, $\alpha$ , $\beta$ , $\gamma$	84.02Å 84.02Å 190.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.40	Depositor
% Data completeness (in resolution range)	92.0 (20.00-2.40)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
Refinement program	TNT	Depositor
R, $R_{free}$	0.191 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4686	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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	1.17	7/1862 (0.4%)	1.77	32/2543 (1.3%)
1	B	1.14	7/1862 (0.4%)	1.80	43/2543 (1.7%)
2	C	1.50	5/457 (1.1%)	2.31	25/615 (4.1%)
2	D	1.45	5/457 (1.1%)	2.09	20/615 (3.3%)
All	All	1.22	24/4638 (0.5%)	1.87	120/6316 (1.9%)

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	B	1	1
2	C	1	0
All	All	2	1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	GLU	CD-OE1	8.25	1.34	1.25
2	D	13	GLU	CD-OE2	8.25	1.34	1.25
2	D	25	GLU	CD-OE1	8.13	1.34	1.25
2	C	25	GLU	CD-OE1	8.06	1.34	1.25
2	D	3	GLU	CD-OE2	7.91	1.34	1.25
1	B	62	GLU	CD-OE2	7.56	1.33	1.25
1	B	70	GLU	CD-OE1	7.19	1.33	1.25
2	C	13	GLU	CD-OE2	7.00	1.33	1.25
2	D	9	GLU	CD-OE1	6.71	1.33	1.25
1	A	62	GLU	CD-OE1	6.27	1.32	1.25
1	A	21	GLU	CD-OE2	6.22	1.32	1.25
1	B	197	GLY	N-CA	6.11	1.55	1.46
1	B	80	GLU	CD-OE1	5.99	1.32	1.25
2	C	18	GLU	CD-OE2	5.82	1.32	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	GLU	CD-OE1	5.81	1.32	1.25
2	D	61	PRO	C-OXT	5.76	1.34	1.23
1	A	80	GLU	CD-OE2	5.74	1.31	1.25
1	B	21	GLU	CD-OE2	5.47	1.31	1.25
1	B	197	GLY	C-O	5.41	1.32	1.23
1	A	183	ALA	N-CA	5.39	1.57	1.46
1	B	43	GLY	C-O	5.19	1.31	1.23
2	C	61	PRO	C-OXT	5.17	1.33	1.23
1	A	193	GLY	C-O	5.16	1.31	1.23
1	A	43	GLY	C-O	5.02	1.31	1.23

All (120) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	48	ARG	NE-CZ-NH1	15.96	128.28	120.30
1	B	65(A)	ARG	NE-CZ-NH2	-15.28	112.66	120.30
1	A	65(A)	ARG	NE-CZ-NH1	14.53	127.56	120.30
1	A	107	ARG	NE-CZ-NH1	13.82	127.21	120.30
2	C	51	ASP	N-CA-CB	-13.38	86.51	110.60
2	D	48	ARG	NE-CZ-NH2	-13.04	113.78	120.30
2	C	24	ASP	CB-CG-OD1	12.80	129.82	118.30
1	B	65(A)	ARG	NE-CZ-NH1	12.67	126.64	120.30
2	C	51	ASP	CB-CG-OD2	-11.59	107.87	118.30
1	A	107	ARG	NE-CZ-NH2	-10.74	114.93	120.30
1	B	98	ASP	CB-CG-OD2	-9.90	109.39	118.30
2	C	34	ARG	NE-CZ-NH1	9.81	125.20	120.30
2	D	35	ARG	NE-CZ-NH1	9.77	125.18	120.30
1	A	194	ASP	CB-CG-OD1	-9.57	109.69	118.30
2	C	24	ASP	CB-CG-OD2	-9.26	109.96	118.30
1	A	60	ASP	CB-CG-OD1	-9.25	109.98	118.30
2	C	34	ARG	NE-CZ-NH2	-9.15	115.72	120.30
2	C	59	GLN	CA-CB-CG	-8.75	94.15	113.40
1	B	60	ASP	CB-CG-OD2	-8.56	110.60	118.30
2	C	47	ARG	NE-CZ-NH2	-8.39	116.10	120.30
1	B	186	ASP	CB-CG-OD1	-8.38	110.76	118.30
1	B	186	ASP	CB-CG-OD2	8.37	125.83	118.30
2	C	35	ARG	NE-CZ-NH1	8.35	124.47	120.30
2	D	47	ARG	NE-CZ-NH1	8.32	124.46	120.30
2	D	44	ARG	NE-CZ-NH2	-8.21	116.19	120.30
1	B	223	ARG	NE-CZ-NH1	8.20	124.40	120.30
1	A	24	ARG	NE-CZ-NH1	-8.11	116.25	120.30
1	B	190	GLY	O-C-N	-8.10	109.73	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	44	ARG	NE-CZ-NH1	8.10	124.35	120.30
1	B	26	SER	N-CA-CB	-8.08	98.38	110.50
1	B	61	ARG	NE-CZ-NH1	7.86	124.23	120.30
1	A	164	ASP	CB-CG-OD2	-7.84	111.25	118.30
1	B	98	ASP	CB-CG-OD1	7.83	125.35	118.30
1	A	216	VAL	CG1-CB-CG2	-7.82	98.39	110.90
2	D	24	ASP	CB-CG-OD1	7.74	125.27	118.30
1	A	164	ASP	CB-CG-OD1	7.71	125.24	118.30
1	B	107	ARG	NE-CZ-NH1	7.71	124.15	120.30
1	B	217(A)	ARG	NE-CZ-NH1	7.70	124.15	120.30
2	D	23	PRO	CA-N-CD	-7.70	100.72	111.50
1	B	90	VAL	CA-CB-CG2	7.69	122.44	110.90
1	A	60	ASP	CB-CG-OD2	7.66	125.19	118.30
1	A	194	ASP	CB-CG-OD2	7.62	125.16	118.30
1	B	217(A)	ARG	NE-CZ-NH2	-7.59	116.50	120.30
2	C	47	ARG	CG-CD-NE	-7.58	95.89	111.80
2	D	1	GLY	O-C-N	7.57	134.82	122.70
2	C	23	PRO	CA-N-CD	-7.53	100.95	111.50
1	B	24	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	B	24	ARG	NE-CZ-NH2	-7.37	116.61	120.30
1	B	162	THR	CA-CB-CG2	7.35	122.69	112.40
2	C	50	ASN	C-N-CA	7.25	139.84	121.70
2	D	24	ASP	CB-CG-OD2	-7.14	111.87	118.30
2	D	32	MET	CG-SD-CE	-7.10	88.84	100.20
2	C	53	LYS	CA-CB-CG	-7.03	97.93	113.40
1	B	171	TYR	CB-CG-CD1	7.01	125.21	121.00
1	A	65	PHE	CB-CA-C	-6.96	96.47	110.40
2	C	51	ASP	CB-CG-OD1	6.85	124.46	118.30
2	C	3	GLU	N-CA-CB	6.84	122.92	110.60
1	A	77	ASP	CB-CG-OD1	-6.83	112.15	118.30
1	B	203	VAL	CB-CA-C	6.71	124.16	111.40
1	B	216	VAL	N-CA-CB	6.71	126.27	111.50
1	B	97	ASP	CB-CG-OD1	-6.71	112.27	118.30
1	A	24	ARG	C-N-CA	-6.71	104.94	121.70
2	C	23	PRO	N-CD-CG	6.65	113.17	103.20
1	B	93	TYR	CB-CG-CD1	-6.58	117.05	121.00
1	B	96	THR	CA-CB-CG2	-6.58	103.19	112.40
1	A	125	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	B	176	VAL	CA-CB-CG2	-6.56	101.05	110.90
1	A	145	ARG	NE-CZ-NH1	6.51	123.56	120.30
2	C	20	LYS	CB-CA-C	-6.43	97.53	110.40
1	A	36	ARG	NE-CZ-NH2	6.27	123.44	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	23	PRO	N-CD-CG	6.27	112.61	103.20
1	B	97	ASP	CB-CG-OD2	6.26	123.94	118.30
1	A	99	VAL	CB-CA-C	-6.24	99.54	111.40
1	B	197	GLY	N-CA-C	-6.19	97.64	113.10
1	A	216	VAL	N-CA-CB	6.16	125.04	111.50
1	B	191	CYS	N-CA-C	6.14	127.57	111.00
1	A	97	ASP	CB-CG-OD2	-6.08	112.83	118.30
2	C	2	GLN	CB-CA-C	6.05	122.50	110.40
2	C	10	VAL	CA-CB-CG2	6.04	119.96	110.90
1	B	164	ASP	CB-CG-OD1	5.99	123.69	118.30
1	A	145	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	B	194	ASP	CB-CG-OD2	-5.91	112.98	118.30
1	B	90	VAL	CG1-CB-CG2	5.91	120.35	110.90
1	B	151	LEU	CB-CG-CD1	-5.89	100.99	111.00
2	D	60	CYS	C-N-CD	-5.83	107.77	120.60
1	B	117	TYR	CB-CG-CD2	-5.83	117.50	121.00
2	D	35	ARG	CD-NE-CZ	5.83	131.75	123.60
2	D	48	ARG	CD-NE-CZ	5.81	131.73	123.60
2	C	15	THR	N-CA-CB	-5.80	99.27	110.30
1	A	48	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	B	194	ASP	CB-CG-OD1	5.75	123.48	118.30
2	D	48	ARG	N-CA-CB	5.75	120.95	110.60
1	B	223	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	35	TYR	N-CA-CB	5.66	120.79	110.60
1	B	207	TYR	CB-CA-C	-5.65	99.09	110.40
1	A	65(A)	ARG	NE-CZ-NH2	-5.60	117.50	120.30
2	D	34	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	A	24	ARG	O-C-N	-5.35	114.14	122.70
1	A	184	GLY	O-C-N	-5.35	114.10	123.20
2	C	2	GLN	N-CA-CB	5.30	120.15	110.60
1	B	45	THR	CA-CB-OG1	5.26	120.05	109.00
1	B	107	ARG	NE-CZ-NH2	-5.25	117.67	120.30
2	C	9	GLU	CG-CD-OE2	-5.24	107.82	118.30
1	A	135	PRO	CB-CA-C	-5.21	98.96	112.00
1	A	61	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	A	97	ASP	CB-CG-OD1	5.21	122.98	118.30
1	B	190	GLY	CA-C-N	5.19	128.63	117.20
1	B	228	PHE	CB-CA-C	-5.15	100.09	110.40
1	A	228	PHE	CB-CA-C	-5.14	100.11	110.40
1	A	186	ASP	CB-CG-OD2	5.14	122.92	118.30
1	B	208	ALA	N-CA-CB	5.12	117.27	110.10
1	A	165	TYR	CB-CA-C	5.12	120.63	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	39	GLU	N-CA-CB	-5.11	101.41	110.60
1	B	61	ARG	CD-NE-CZ	5.09	130.73	123.60
2	D	24	ASP	N-CA-CB	-5.05	101.52	110.60
2	D	1	GLY	CA-C-N	-5.04	106.11	117.20
1	B	60	ASP	CB-CG-OD1	5.03	122.82	118.30
2	D	1	GLY	C-N-CA	5.01	134.22	121.70
2	C	1	GLY	N-CA-C	-5.00	100.59	113.10
2	C	21	CYS	C-N-CA	-5.00	111.79	122.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	203	VAL	CA
2	C	2	GLN	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	30	GLN	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	1822	0	1757	20	0
1	B	1822	0	1757	64	0
2	C	448	0	408	12	0
2	D	448	0	409	10	0
3	A	58	0	0	0	0
3	B	61	0	0	5	0
3	C	16	0	0	1	0
3	D	11	0	0	0	0
All	All	4686	0	4331	103	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 12.

All (103) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:VAL:HG21	1:B:156:GLN:HB2	1.33	1.08
1:B:115:ASN:ND2	1:B:117:TYR:H	1.52	1.07
1:B:61:ARG:HB3	1:B:63:LEU:HD13	1.36	1.03
2:C:24:ASP:OD1	2:C:24:ASP:N	1.97	0.97
1:B:125:ARG:HG3	1:B:125:ARG:HH11	1.34	0.92
1:B:16:VAL:HG11	1:B:156:GLN:HB3	1.53	0.90
1:B:115:ASN:HD22	1:B:117:TYR:H	1.16	0.88
1:B:61:ARG:HB3	1:B:63:LEU:CD1	2.03	0.87
2:D:18:GLU:OE1	2:D:48:ARG:HD2	1.82	0.79
1:B:50:ASN:HD22	1:B:50:ASN:H	1.28	0.78
1:B:62:GLU:C	1:B:63:LEU:HD12	2.04	0.77
1:B:16:VAL:HG11	1:B:156:GLN:CB	2.13	0.77
1:B:23:GLN:O	1:B:26:SER:HB2	1.86	0.76
2:C:35:ARG:HG2	3:C:74:HOH:O	1.86	0.74
1:B:115:ASN:HD22	1:B:116:SER:N	1.85	0.74
2:D:24:ASP:OD1	2:D:24:ASP:N	2.17	0.74
1:A:36:ARG:NH1	2:D:51:ASP:O	2.23	0.72
2:D:20:LYS:O	2:D:23:PRO:HD2	1.89	0.72
1:B:125:ARG:HG3	1:B:125:ARG:NH1	2.02	0.71
1:B:16:VAL:CG2	1:B:156:GLN:HB2	2.17	0.70
2:C:2:GLN:HG3	2:C:11:TRP:CG	2.26	0.70
1:B:115:ASN:HD21	1:B:117:TYR:HD2	1.39	0.69
1:A:187:GLY:HA3	1:A:222:THR:HG22	1.76	0.67
1:B:33:LEU:O	1:B:41:THR:HG22	1.95	0.67
1:B:115:ASN:HD22	1:B:117:TYR:N	1.93	0.65
2:C:20:LYS:O	2:C:23:PRO:HD2	1.98	0.63
1:B:16:VAL:HG12	3:B:271:HOH:O	1.98	0.62
1:B:197:GLY:HA2	1:B:213:THR:H	1.63	0.62
2:C:49:THR:OG1	2:C:51:ASP:HB3	1.99	0.61
1:B:60:ASP:OD2	1:B:96:THR:HG23	2.01	0.61
1:B:50:ASN:N	1:B:50:ASN:HD22	1.94	0.60
2:C:1:GLY:HA3	2:C:11:TRP:H	1.67	0.59
1:B:133:ASN:ND2	1:B:133:ASN:O	2.30	0.59
1:B:16:VAL:HB	3:B:272:HOH:O	2.01	0.59
1:B:62:GLU:CD	1:B:62:GLU:H	2.06	0.59
1:B:75:GLN:O	3:B:246:HOH:O	2.17	0.59
1:A:216:VAL:HG13	2:C:31:LEU:HD11	1.85	0.58
1:B:62:GLU:N	1:B:62:GLU:OE1	2.35	0.58
1:A:38:TRP:CE3	1:A:65(A):ARG:HD2	2.39	0.58
1:B:218:LEU:H	1:B:218:LEU:HD12	1.68	0.57
1:B:91:HIS:CG	1:B:92:PRO:HD2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:20:LYS:C	2:D:23:PRO:HD2	2.27	0.55
2:C:23:PRO:HA	2:C:24:ASP:OD1	2.06	0.55
1:B:134:SER:H	1:B:162:THR:HB	1.73	0.54
2:D:49:THR:OG1	2:D:53:LYS:HG3	2.07	0.54
1:B:215:PHE:CD2	2:D:28:PRO:HB3	2.43	0.54
1:B:197:GLY:CA	1:B:213:THR:H	2.22	0.53
1:B:221(A):VAL:HG22	1:B:224:LYS:HB2	1.91	0.53
1:B:62:GLU:O	1:B:63:LEU:HD12	2.09	0.53
1:B:244:SER:OG	1:B:244:SER:O	2.27	0.53
2:C:2:GLN:HG3	2:C:11:TRP:CB	2.39	0.52
1:B:115:ASN:ND2	1:B:117:TYR:N	2.37	0.52
2:C:55:ILE:HG13	2:C:55:ILE:O	2.09	0.52
1:B:133:ASN:HD22	1:B:133:ASN:C	2.11	0.51
1:B:125:ARG:HH11	1:B:125:ARG:CG	2.17	0.51
1:B:131:ALA:O	1:B:162:THR:HG21	2.11	0.51
1:B:81:GLN:HE22	1:B:113:THR:H	1.57	0.51
1:B:188:VAL:HG12	1:B:188(A):ARG:HG3	1.93	0.51
1:B:115:ASN:ND2	1:B:117:TYR:CD2	2.78	0.51
1:B:82:TYR:CD1	1:B:82:TYR:N	2.79	0.50
1:B:68:VAL:CG2	1:B:81:GLN:HB2	2.42	0.50
1:B:149:GLY:HA3	3:B:282:HOH:O	2.11	0.50
1:B:223:ARG:NH1	3:B:277:HOH:O	2.28	0.49
1:B:144:THR:O	1:B:145:ARG:HB3	2.12	0.49
1:A:90:VAL:HG12	1:A:91:HIS:N	2.28	0.48
1:B:37:SER:HB2	1:B:38:TRP:H	1.45	0.48
1:B:115:ASN:C	1:B:115:ASN:HD22	2.17	0.48
1:B:68:VAL:HG23	1:B:81:GLN:HB2	1.96	0.47
1:B:134:SER:N	1:B:162:THR:HB	2.30	0.47
1:B:81:GLN:NE2	1:B:112:VAL:HG13	2.29	0.47
2:D:1:GLY:N	2:D:5:CYS:HB2	2.30	0.47
1:A:90:VAL:CG1	1:A:91:HIS:N	2.78	0.46
2:D:36:PRO:O	2:D:37:SER:HB3	2.16	0.46
1:B:61:ARG:HE	1:B:63:LEU:HD11	1.80	0.46
1:B:50:ASN:N	1:B:50:ASN:ND2	2.61	0.46
1:B:44:GLY:HA2	1:B:196:GLY:O	2.16	0.46
1:B:217:SER:OG	1:B:218:LEU:HD12	2.15	0.46
1:B:221:ASN:OD1	1:B:221:ASN:N	2.40	0.46
1:A:68:VAL:HG22	1:A:81:GLN:HB2	1.98	0.45
1:A:109:ALA:C	1:A:110:GLN:HG3	2.35	0.45
1:B:16:VAL:O	1:B:16:VAL:HG12	2.15	0.45
2:D:41:SER:HA	2:D:42:PRO:HD3	1.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:GLY:O	1:B:194:ASP:OD2	2.34	0.45
1:A:172:TRP:CD1	1:A:225:PRO:HD2	2.52	0.44
1:A:68:VAL:CG2	1:A:81:GLN:HB2	2.46	0.44
1:A:240:ASN:OD1	1:A:240:ASN:N	2.49	0.43
2:C:49:THR:OG1	2:C:53:LYS:HG3	2.19	0.43
1:A:115:ASN:OD1	1:A:116:SER:N	2.52	0.42
1:B:86:GLN:NE2	1:B:109:ALA:HA	2.34	0.42
1:A:82:TYR:CD1	1:A:82:TYR:N	2.86	0.42
1:B:31:ILE:HD12	1:B:31:ILE:C	2.40	0.42
1:A:37:SER:HB2	1:A:38:TRP:H	1.68	0.41
1:B:216:VAL:HB	1:B:217:SER:H	1.74	0.41
1:A:27:TRP:N	1:A:28:PRO:CD	2.82	0.41
1:B:17:VAL:HG23	1:B:191:CYS:HB2	2.01	0.41
2:C:30:PRO:HD2	2:C:34:ARG:NH2	2.36	0.41
1:A:202:LEU:HD22	1:A:207:TYR:CE1	2.55	0.41
1:A:172:TRP:NE1	1:A:225:PRO:HD2	2.36	0.41
1:A:28:PRO:HB2	1:A:119:GLN:HB2	2.02	0.41
1:A:100:GLY:O	1:A:101:TYR:HB2	2.22	0.40
1:A:90:VAL:HG12	1:A:91:HIS:O	2.21	0.40
1:B:99:VAL:H	1:B:99:VAL:HG23	1.48	0.40
1:B:62:GLU:O	1:B:62:GLU:HG2	2.21	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	238/240 (99%)	221 (93%)	16 (7%)	1 (0%)	39	56
1	B	238/240 (99%)	220 (92%)	18 (8%)	0	100	100
2	C	59/61 (97%)	56 (95%)	3 (5%)	0	100	100
2	D	59/61 (97%)	53 (90%)	4 (7%)	2 (3%)	5	4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	594/602 (99%)	550 (93%)	41 (7%)	3 (0%)	34 48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	2	GLN
2	D	3	GLU
1	A	28	PRO

### 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	198/198 (100%)	182 (92%)	16 (8%)	15 22
1	B	198/198 (100%)	178 (90%)	20 (10%)	9 13
2	C	53/53 (100%)	49 (92%)	4 (8%)	17 26
2	D	53/53 (100%)	50 (94%)	3 (6%)	25 40
All	All	502/502 (100%)	459 (91%)	43 (9%)	13 19

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36(C)	SER
1	A	60	ASP
1	A	61	ARG
1	A	65(A)	ARG
1	A	77	ASP
1	A	86	GLN
1	A	96	THR
1	A	97	ASP
1	A	99	VAL
1	A	105	LEU
1	A	143	LEU
1	A	181	VAL

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Mol	Chain	Res	Type
1	A	188(A)	ARG
1	A	195	SER
1	A	203	VAL
1	A	222	THR
1	B	26	SER
1	B	36(C)	SER
1	B	45	THR
1	B	50	ASN
1	B	76	ASN
1	B	86	GLN
1	B	90	VAL
1	B	96	THR
1	B	115	ASN
1	B	125	ARG
1	B	130	LEU
1	B	133	ASN
1	B	153	GLN
1	B	170(A)	SER
1	B	170(B)	SER
1	B	178	ASN
1	B	191	CYS
1	B	202	LEU
1	B	221(A)	VAL
1	B	231	VAL
2	C	2	GLN
2	C	24	ASP
2	C	41	SER
2	C	53	LYS
2	D	24	ASP
2	D	35	ARG
2	D	41	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	A	119	GLN
1	A	239	ASN
1	B	50	ASN
1	B	76	ASN
1	B	81	GLN
1	B	115	ASN

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Mol	Chain	Res	Type
1	B	204	ASN

### 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 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates ⓘ

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

### 6.4 Ligands ⓘ

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

### 6.5 Other polymers ⓘ

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