



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:19 AM UTC

PDB ID : 1FLE / pdb_00001fle
Title : CRYSTAL STRUCTURE OF ELAFIN COMPLEXED WITH PORCINE PANCREATIC ELASTASE
Authors : Tsunemi, M.; Matsuura, Y.; Sakakibara, S.; Katsube, Y.
Deposited on : 1996-07-04
Resolution : 1.90 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
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

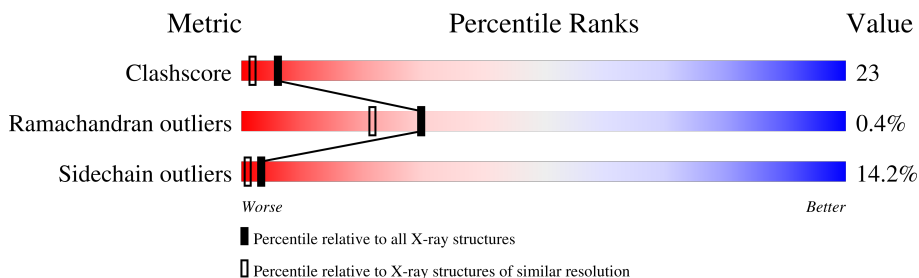
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	240	
2	I	57	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2320 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 ELASTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	240	1822	1135	330	347	10	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	77	ASN	ASP	conflict	UNP P00772

- Molecule 2 is a protein called ELAFIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	I	47	341	210	60	61	10	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	142	Total	O	0	0
			142	142		
3	I	15	Total	O	0	0
			15	15		

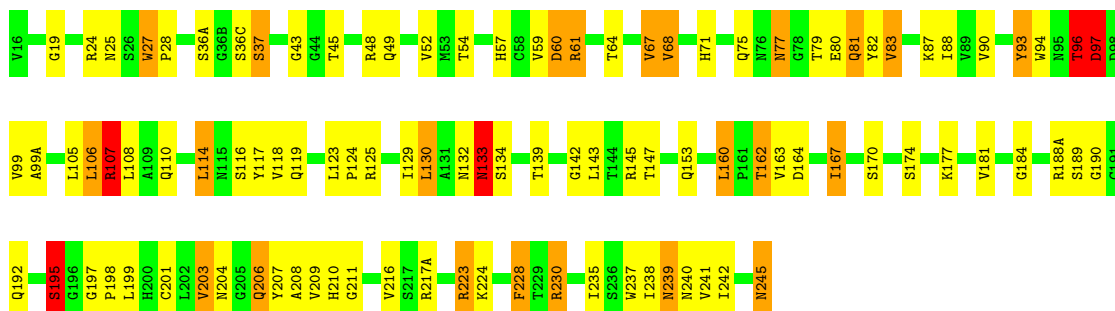
3 Residue-property plots [i](#)

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. 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. 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.

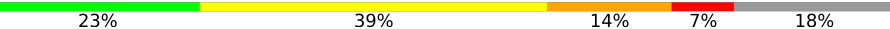
Note EDS was not executed.

- Molecule 1: ELASTASE

Chain E:  57% 31% 10%



- Molecule 2: ELAFIN

Chain I:  23% 39% 14% 7% 18%



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	37.91Å 73.32Å 48.92Å 90.00° 105.40° 90.00°	Depositor
Resolution (Å)	10.00 – 1.90	Depositor
% Data completeness (in resolution range)	76.8 (10.00-1.90)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.197 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	2320	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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	E	1.33	5/1862 (0.3%)	2.23	69/2543 (2.7%)
2	I	1.22	0/347	2.43	19/466 (4.1%)
All	All	1.31	5/2209 (0.2%)	2.26	88/3009 (2.9%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	199	LEU	CA-C	6.00	1.60	1.52
1	E	230	ARG	N-CA	5.38	1.53	1.46
1	E	54	THR	CA-CB	5.37	1.61	1.53
1	E	197	GLY	N-CA	5.32	1.51	1.44
1	E	211	GLY	N-CA	5.12	1.50	1.45

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	48	ARG	CD-NE-CZ	39.56	179.79	124.40
2	I	41	ILE	CA-C-O	14.60	139.03	120.78
2	I	22	ARG	CD-NE-CZ	13.75	143.65	124.40
1	E	240	ASN	CA-CB-CG	12.31	124.91	112.60
1	E	48	ARG	NE-CZ-NH2	11.15	129.23	119.20
1	E	164	ASP	CA-CB-CG	11.03	123.62	112.60
1	E	133	ASN	OD1-CG-ND2	9.78	132.38	122.60
1	E	125	ARG	NE-CZ-NH2	9.49	127.74	119.20
1	E	107	ARG	CD-NE-CZ	8.08	135.71	124.40
2	I	35	ASP	CA-CB-CG	8.05	120.65	112.60
1	E	190	GLY	N-CA-C	-7.98	100.00	112.85
2	I	42	LYS	CA-C-N	-7.87	111.69	122.86
2	I	42	LYS	C-N-CA	-7.87	111.69	122.86
1	E	125	ARG	CD-NE-CZ	7.76	135.27	124.40
2	I	19	ILE	CB-CA-C	7.76	121.79	110.77
2	I	28	PRO	O-C-N	7.71	124.77	121.15
1	E	208	ALA	O-C-N	7.70	132.03	123.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	77	ASN	CA-CB-CG	7.57	120.17	112.60
1	E	206	GLN	CB-CG-CD	7.57	125.46	112.60
2	I	41	ILE	CB-CA-C	7.32	123.29	111.29
1	E	108	LEU	O-C-N	7.27	131.18	122.96
1	E	36(C)	SER	N-CA-C	-7.22	105.39	114.56
1	E	68	VAL	O-C-N	7.17	130.44	122.98
1	E	64	THR	N-CA-CB	7.09	121.97	110.42
2	I	18	ILE	N-CA-CB	7.03	119.89	110.26
2	I	46	GLU	CB-CG-CD	6.96	124.44	112.60
1	E	195	SER	CB-CA-C	-6.84	96.80	110.42
1	E	97	ASP	N-CA-CB	-6.80	100.12	110.26
1	E	43	GLY	N-CA-C	-6.76	102.25	112.60
2	I	12	LYS	O-C-N	6.70	128.29	121.72
1	E	223	ARG	CD-NE-CZ	-6.58	115.18	124.40
1	E	199	LEU	N-CA-CB	6.56	121.02	110.65
1	E	217(A)	ARG	CD-NE-CZ	-6.53	115.26	124.40
1	E	203	VAL	N-CA-CB	-6.48	102.40	111.99
1	E	36(C)	SER	O-C-N	6.48	128.97	122.10
1	E	162	THR	CA-CB-OG1	-6.41	99.98	109.60
1	E	61	ARG	CD-NE-CZ	6.40	133.35	124.40
1	E	237	TRP	CA-C-N	6.36	128.81	120.60
1	E	237	TRP	C-N-CA	6.36	128.81	120.60
2	I	57	GLN	OE1-CD-NE2	6.36	128.96	122.60
2	I	22	ARG	NE-CZ-NH1	6.32	127.82	121.50
2	I	35	ASP	N-CA-C	-6.24	104.53	111.71
1	E	60	ASP	CA-CB-CG	6.13	118.73	112.60
1	E	163	VAL	N-CA-C	-5.98	97.77	107.28
2	I	44	CYS	N-CA-C	-5.96	100.13	109.25
1	E	59	VAL	O-C-N	5.93	128.56	122.09
1	E	199	LEU	CA-C-O	-5.87	114.24	120.40
1	E	245	ASN	CA-CB-CG	5.82	118.42	112.60
2	I	43	LYS	N-CA-CB	5.79	119.64	110.84
1	E	81	GLN	OE1-CD-NE2	5.74	128.34	122.60
1	E	170	SER	N-CA-CB	5.72	118.54	110.36
1	E	25	ASN	CA-CB-CG	5.71	118.31	112.60
1	E	125	ARG	NH1-CZ-NH2	-5.70	111.89	119.30
1	E	27	TRP	CA-C-N	5.69	125.40	119.82
1	E	27	TRP	C-N-CA	5.69	125.40	119.82
1	E	67	VAL	O-C-N	5.63	129.34	123.26
1	E	203	VAL	CB-CA-C	5.45	117.53	110.12
1	E	71	HIS	N-CA-CB	5.44	119.68	110.49
1	E	240	ASN	N-CA-CB	5.34	118.07	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	94	TRP	CA-C-N	5.32	131.01	123.14
1	E	94	TRP	C-N-CA	5.32	131.01	123.14
2	I	52	ALA	CA-C-N	5.32	129.49	122.42
2	I	52	ALA	C-N-CA	5.32	129.49	122.42
1	E	134	SER	O-C-N	5.31	127.37	121.53
1	E	48	ARG	NE-CZ-NH1	-5.29	116.20	121.50
1	E	238	ILE	O-C-N	5.28	127.22	121.94
1	E	224	LYS	O-C-N	5.27	126.11	121.37
1	E	93	TYR	O-C-N	5.26	128.92	122.34
1	E	145	ARG	CD-NE-CZ	5.25	131.75	124.40
1	E	97	ASP	N-CA-C	5.23	119.72	112.45
1	E	216	VAL	O-C-N	5.22	128.28	122.75
1	E	133	ASN	CA-CB-CG	-5.19	107.41	112.60
1	E	204	ASN	CA-C-O	-5.18	115.53	121.54
1	E	170	SER	N-CA-C	-5.16	103.72	110.53
1	E	147	THR	N-CA-CB	5.14	117.44	109.48
1	E	83	VAL	O-C-N	5.14	128.75	123.05
1	E	133	ASN	O-C-N	5.13	130.77	122.20
1	E	170	SER	CA-CB-OG	5.12	121.35	111.10
1	E	167	ILE	O-C-N	5.11	128.54	121.90
1	E	133	ASN	CA-C-O	-5.10	114.45	120.57
1	E	64	THR	O-C-N	5.09	129.02	123.22
1	E	198	PRO	N-CA-C	5.08	119.03	111.41
1	E	199	LEU	N-CA-C	-5.08	99.13	108.02
1	E	217(A)	ARG	N-CA-CB	5.05	118.08	110.30
1	E	80	GLU	N-CA-C	5.05	117.95	110.48
1	E	19	GLY	CA-C-N	5.03	130.46	122.62
1	E	19	GLY	C-N-CA	5.03	130.46	122.62
2	I	23	CYS	N-CA-C	-5.01	103.53	110.35

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	E	1822	0	1759	51	2
2	I	341	0	344	52	2
3	E	142	0	0	13	4
3	I	15	0	0	0	0
All	All	2320	0	2103	97	4

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 (97) 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:E:203:VAL:HB	3:E:541:HOH:O	1.29	1.26
2:I:12:LYS:HG2	2:I:43:LYS:NZ	1.53	1.20
2:I:12:LYS:HG2	2:I:43:LYS:HZ3	1.05	1.01
2:I:38:CYS:O	2:I:42:LYS:HB3	1.60	1.01
2:I:35:ASP:O	2:I:42:LYS:HB2	1.60	1.00
1:E:189:SER:N	3:E:512:HOH:O	1.95	0.99
2:I:12:LYS:CG	2:I:43:LYS:HZ3	1.73	0.99
2:I:12:LYS:HB3	2:I:35:ASP:OD2	1.63	0.98
2:I:12:LYS:HE3	2:I:14:GLY:O	1.62	0.97
2:I:22:ARG:NH2	2:I:28:PRO:HG3	1.79	0.96
1:E:97:ASP:OD2	2:I:31:ARG:NH2	1.99	0.95
2:I:12:LYS:HE3	2:I:43:LYS:HE2	1.52	0.90
1:E:192:GLN:HB3	2:I:26:LEU:HD13	1.53	0.89
2:I:22:ARG:HH22	2:I:28:PRO:HG3	1.39	0.86
1:E:167:ILE:HD13	3:E:503:HOH:O	1.76	0.84
2:I:12:LYS:CE	2:I:43:LYS:HZ3	1.90	0.83
2:I:12:LYS:CE	2:I:14:GLY:O	2.30	0.80
1:E:60:ASP:OD2	1:E:96:THR:HG23	1.82	0.79
2:I:12:LYS:CB	2:I:35:ASP:OD2	2.30	0.78
2:I:12:LYS:HG2	2:I:43:LYS:HZ2	1.45	0.76
2:I:12:LYS:O	2:I:12:LYS:HD3	1.86	0.76
2:I:12:LYS:CG	2:I:43:LYS:NZ	2.39	0.75
2:I:12:LYS:HE3	2:I:43:LYS:CE	2.20	0.71
1:E:49:GLN:NE2	1:E:114:LEU:HD21	2.05	0.71
2:I:12:LYS:CG	2:I:35:ASP:OD2	2.39	0.71
2:I:43:LYS:HG3	2:I:56:PRO:HB3	1.71	0.71
2:I:12:LYS:CD	2:I:43:LYS:HZ3	2.04	0.70
1:E:88:ILE:HG12	1:E:106:LEU:HD12	1.74	0.70
1:E:181:VAL:HG23	1:E:230:ARG:HB2	1.73	0.70
2:I:12:LYS:HA	2:I:42:LYS:HZ1	1.57	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:20:LEU:HD23	2:I:20:LEU:O	1.92	0.69
1:E:97:ASP:O	3:E:437:HOH:O	2.12	0.68
2:I:12:LYS:HA	2:I:42:LYS:NZ	2.10	0.66
1:E:130:LEU:HD11	1:E:210:HIS:CE1	2.33	0.63
1:E:133:ASN:ND2	3:E:485:HOH:O	2.32	0.63
1:E:203:VAL:HG23	3:E:478:HOH:O	1.99	0.62
2:I:46:GLU:H	2:I:46:GLU:CD	2.09	0.60
2:I:35:ASP:HB3	2:I:42:LYS:HE2	1.82	0.60
2:I:12:LYS:CE	2:I:43:LYS:NZ	2.62	0.60
1:E:87:LYS:HB2	1:E:107:ARG:HB2	1.83	0.59
2:I:40:GLY:O	2:I:42:LYS:HG3	2.03	0.59
2:I:17:PRO:HG2	2:I:19:ILE:HD11	1.85	0.59
1:E:105:LEU:HD12	1:E:241:VAL:HG11	1.85	0.58
2:I:12:LYS:NZ	2:I:43:LYS:NZ	2.52	0.58
1:E:132:ASN:O	1:E:133:ASN:HB2	2.04	0.58
1:E:107:ARG:HH21	1:E:245:ASN:C	2.11	0.57
1:E:24:ARG:HH22	1:E:77:ASN:ND2	2.02	0.57
2:I:12:LYS:NZ	2:I:43:LYS:HZ3	2.02	0.57
2:I:45:CYS:HB2	2:I:54:PHE:CD1	2.39	0.57
1:E:68:VAL:CG2	1:E:81:GLN:HB2	2.36	0.55
1:E:162:THR:N	3:E:506:HOH:O	2.40	0.55
1:E:28:PRO:HB2	1:E:119:GLN:HB2	1.88	0.55
2:I:12:LYS:HG2	2:I:35:ASP:OD2	2.08	0.54
1:E:79:THR:HG22	1:E:117:TYR:CE2	2.42	0.54
1:E:203:VAL:CB	3:E:541:HOH:O	2.13	0.54
2:I:14:GLY:O	2:I:43:LYS:HE2	2.07	0.54
1:E:192:GLN:CB	2:I:26:LEU:HD13	2.31	0.53
1:E:96:THR:HG22	3:E:431:HOH:O	2.07	0.53
1:E:68:VAL:HG22	1:E:81:GLN:HB2	1.90	0.53
1:E:160:LEU:HD22	1:E:184:GLY:HA3	1.89	0.53
2:I:34:LYS:O	2:I:37:ASP:HB2	2.10	0.51
1:E:45:THR:HG21	1:E:209:VAL:HG21	1.93	0.51
2:I:38:CYS:C	2:I:42:LYS:HB3	2.34	0.49
1:E:36(A):SER:O	1:E:37:SER:OG	2.26	0.48
1:E:67:VAL:HG22	1:E:82:TYR:CD2	2.47	0.48
1:E:49:GLN:NE2	1:E:114:LEU:CD2	2.77	0.48
2:I:12:LYS:CB	2:I:43:LYS:HD2	2.44	0.47
1:E:81:GLN:NE2	1:E:118:VAL:HG21	2.30	0.46
1:E:57:HIS:CE1	1:E:195:SER:HG	2.26	0.46
1:E:99(A):ALA:HB2	2:I:21:ILE:CD1	2.46	0.46
1:E:162:THR:CA	3:E:506:HOH:O	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:12:LYS:HB3	2:I:43:LYS:HD2	1.98	0.45
2:I:22:ARG:HH22	2:I:28:PRO:CG	2.20	0.45
1:E:87:LYS:O	1:E:106:LEU:HA	2.17	0.45
1:E:129:ILE:HD12	3:E:461:HOH:O	2.16	0.45
1:E:201:CYS:O	1:E:207:TYR:HA	2.16	0.44
2:I:15:SER:O	2:I:56:PRO:HA	2.18	0.44
2:I:34:LYS:C	2:I:36:THR:N	2.72	0.44
1:E:142:GLY:O	1:E:143:LEU:C	2.61	0.44
2:I:17:PRO:HG2	2:I:19:ILE:CD1	2.48	0.43
2:I:42:LYS:N	2:I:42:LYS:HD3	2.34	0.43
2:I:12:LYS:O	2:I:12:LYS:CD	2.63	0.43
1:E:235:ILE:O	1:E:239:ASN:HB2	2.19	0.43
2:I:35:ASP:O	2:I:42:LYS:CB	2.49	0.43
2:I:12:LYS:HB2	2:I:43:LYS:CD	2.49	0.43
1:E:160:LEU:HD22	1:E:184:GLY:CA	2.49	0.42
1:E:99(A):ALA:HB2	2:I:21:ILE:HD13	2.02	0.42
1:E:167:ILE:CD1	3:E:503:HOH:O	2.53	0.42
1:E:189:SER:CA	3:E:512:HOH:O	2.57	0.42
1:E:192:GLN:HA	2:I:24:ALA:C	2.45	0.41
1:E:223:ARG:HH11	1:E:223:ARG:HD2	1.55	0.41
1:E:123:LEU:HA	1:E:124:PRO:HD3	1.96	0.41
1:E:27:TRP:CD1	1:E:139:THR:HG21	2.56	0.41
1:E:228:PHE:N	1:E:228:PHE:CD1	2.90	0.40
1:E:93:TYR:N	1:E:93:TYR:CD1	2.89	0.40
1:E:188(A):ARG:O	1:E:189:SER:HB3	2.22	0.40
2:I:43:LYS:O	2:I:53:CYS:HA	2.21	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:13:PRO:O	3:E:528:HOH:O[2_656]	1.79	0.41
1:E:153:GLN:CD	3:E:492:HOH:O[2_657]	2.14	0.06
2:I:12:LYS:NZ	3:E:527:HOH:O[2_656]	2.16	0.04
1:E:153:GLN:CG	3:E:492:HOH:O[2_657]	2.18	0.02

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	E	238/240 (99%)	226 (95%)	11 (5%)	1 (0%)	30	22
2	I	45/57 (79%)	40 (89%)	5 (11%)	0	100	100
All	All	283/297 (95%)	266 (94%)	16 (6%)	1 (0%)	30	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	96	THR

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	E	198/198 (100%)	174 (88%)	24 (12%)	5	2
2	I	41/49 (84%)	31 (76%)	10 (24%)	1	0
All	All	239/247 (97%)	205 (86%)	34 (14%)	3	1

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	37	SER
1	E	52	VAL
1	E	61	ARG
1	E	75	GLN
1	E	83	VAL

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Mol	Chain	Res	Type
1	E	90	VAL
1	E	96	THR
1	E	97	ASP
1	E	99	VAL
1	E	106	LEU
1	E	107	ARG
1	E	110	GLN
1	E	114	LEU
1	E	116	SER
1	E	130	LEU
1	E	133	ASN
1	E	160	LEU
1	E	174	SER
1	E	177	LYS
1	E	195	SER
1	E	206	GLN
1	E	228	PHE
1	E	239	ASN
1	E	242	ILE
2	I	11	THR
2	I	12	LYS
2	I	20	LEU
2	I	26	LEU
2	I	33	LEU
2	I	36	THR
2	I	41	ILE
2	I	42	LYS
2	I	43	LYS
2	I	46	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	23	GLN
1	E	49	GLN
1	E	119	GLN
1	E	133	ASN
1	E	239	ASN
1	E	240	ASN
2	I	57	GLN

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](#)

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.