



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 08:36 PM EDT

PDB ID : 1LI1
Title : The 1.9-Å crystal structure of the noncollagenous (NC1) domain of human placenta collagen IV shows stabilization via a novel type of covalent Met-Lys cross-link
Authors : Than, M.E.; Henrich, S.; Huber, R.; Ries, A.; Mann, K.; Kuhn, K.; Timpl, R.; Bourenkov, G.P.; Bartunik, H.D.; Bode, W.
Deposited on : 2002-04-17
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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

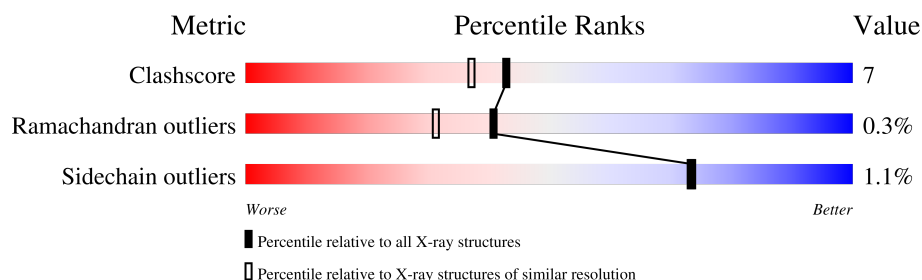
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (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	A	229	84% 15% .
1	B	229	86% 13% .
1	D	229	87% 12%
1	E	229	87% 12% .
2	C	228	82% 16% .
2	F	228	79% 20% .

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	A	2009	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11664 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 Collagen alpha 1(IV).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	24	0	0
			1765	1109	309	326	21			
1	B	227	Total	C	N	O	S	21	0	0
			1758	1104	308	325	21			
1	D	228	Total	C	N	O	S	34	0	0
			1765	1109	309	326	21			
1	E	227	Total	C	N	O	S	22	0	0
			1758	1104	308	325	21			

- Molecule 2 is a protein called Collagen alpha 2(IV).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	224	Total	C	N	O	S	14	0	0
			1730	1096	291	324	19			
2	F	225	Total	C	N	O	S	29	0	0
			1738	1102	292	325	19			

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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	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	D	1	Total C O 4 2 2	0	0
3	E	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	222	Total O 222 222	0	0
4	B	185	Total O 185 185	0	0
4	C	167	Total O 167 167	0	0

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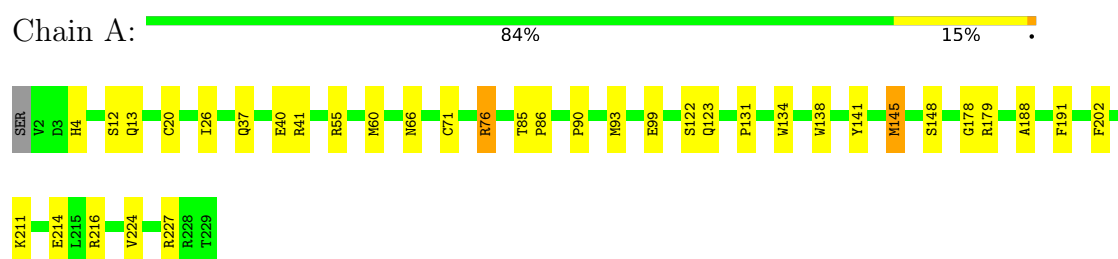
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	208	Total 208	O 208	0	0
4	E	187	Total 187	O 187	0	0
4	F	145	Total 145	O 145	0	0

3 Residue-property plots

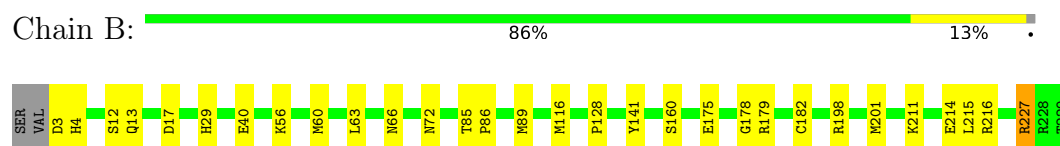
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

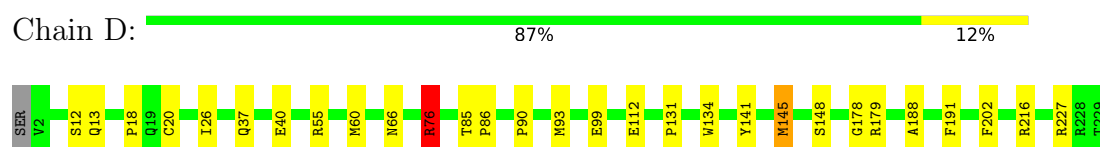
- Molecule 1: Collagen alpha 1(IV)



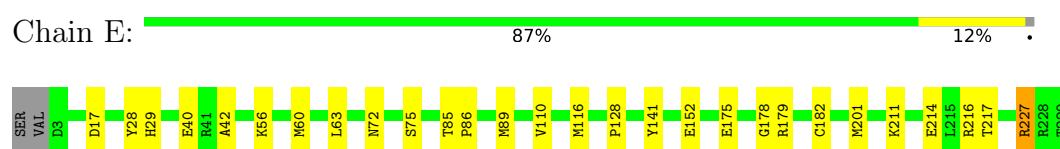
- Molecule 1: Collagen alpha 1(IV)



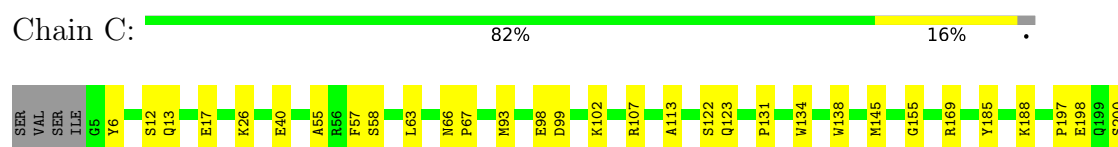
- Molecule 1: Collagen alpha 1(IV)



- Molecule 1: Collagen alpha 1(IV)

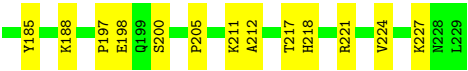
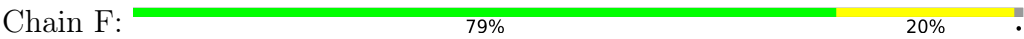


- Molecule 2: Collagen alpha 2(IV)





● Molecule 2: Collagen alpha 2(IV)



4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	234.59Å 234.59Å 99.49Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.95 – 1.90	Depositor
% Data completeness (in resolution range)	92.3 (19.95-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.182 , 0.202	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11664	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.48	0/1816	0.71	0/2465
1	B	0.45	0/1809	0.73	2/2455 (0.1%)
1	D	0.47	0/1816	0.72	1/2465 (0.0%)
1	E	0.47	0/1809	0.73	2/2455 (0.1%)
2	C	0.43	0/1781	0.71	2/2420 (0.1%)
2	F	0.44	0/1789	0.71	2/2431 (0.1%)
All	All	0.46	0/10820	0.72	9/14691 (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
2	F	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	227	ARG	NE-CZ-NH2	-9.02	115.79	120.30
1	E	227	ARG	NE-CZ-NH1	-8.85	115.88	120.30
1	B	227	ARG	NE-CZ-NH1	8.28	124.44	120.30
2	F	169	ARG	NE-CZ-NH1	-7.21	116.69	120.30
2	C	169	ARG	NE-CZ-NH2	-6.82	116.89	120.30
2	F	169	ARG	NE-CZ-NH2	6.58	123.59	120.30
2	C	169	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	D	76	ARG	NE-CZ-NH1	6.01	123.30	120.30
1	E	227	ARG	NE-CZ-NH2	5.71	123.16	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	35	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	1765	0	1680	29	0
1	B	1758	0	1671	25	0
1	D	1765	0	1680	25	0
1	E	1758	0	1671	25	0
2	C	1730	0	1641	26	0
2	F	1738	0	1652	35	0
3	A	4	0	3	2	0
3	B	16	0	12	0	0
3	C	8	0	6	0	0
3	D	4	0	3	1	0
3	E	4	0	3	0	0
4	A	222	0	0	6	0
4	B	185	0	0	6	0
4	C	167	0	0	1	0
4	D	208	0	0	2	0
4	E	187	0	0	6	0
4	F	145	0	0	3	0
All	All	11664	0	10022	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 7.

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)
2:C:93:MET:SD	2:F:211:LYS:HE2	2.01	1.00
2:C:211:LYS:HE3	2:F:93:MET:SD	2.11	0.91
1:A:40:GLU:OE2	1:D:37:GLN:NE2	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:185:TYR:H	2:F:188:LYS:HZ2	1.42	0.67
1:E:72:ASN:HD21	1:E:75:SER:CB	2.12	0.61
1:D:179:ARG:HH11	1:D:179:ARG:HG2	1.66	0.60
2:C:185:TYR:H	2:C:188:LYS:HZ2	1.50	0.59
2:F:107:ARG:HD3	4:F:358:HOH:O	2.01	0.59
1:D:179:ARG:HG2	1:D:179:ARG:NH1	2.17	0.59
1:B:211:LYS:HE3	1:D:93:MET:SD	2.43	0.57
1:E:72:ASN:HD21	1:E:75:SER:HB2	1.70	0.57
2:F:116:ILE:HG13	2:F:117:ALA:N	2.18	0.56
2:F:217:THR:HG23	2:F:218:HIS:ND1	2.21	0.56
1:E:72:ASN:HB2	4:E:2139:HOH:O	2.05	0.56
1:A:211:LYS:O	1:A:214:GLU:HG2	2.07	0.55
2:F:185:TYR:H	2:F:188:LYS:NZ	2.04	0.55
1:A:93:MET:SD	1:E:211:LYS:HE3	2.46	0.54
1:A:179:ARG:HG2	1:A:179:ARG:NH1	2.22	0.54
4:A:2129:HOH:O	1:D:93:MET:HE1	2.07	0.54
1:B:211:LYS:O	1:B:214:GLU:HG2	2.08	0.54
1:E:217:THR:HG23	4:E:2157:HOH:O	2.07	0.54
1:A:99:GLU:HG2	4:A:2203:HOH:O	2.07	0.54
1:A:179:ARG:HG2	1:A:179:ARG:HH11	1.71	0.54
2:C:185:TYR:H	2:C:188:LYS:NZ	2.06	0.53
1:A:37:GLN:NE2	1:D:40:GLU:OE1	2.41	0.53
2:C:63:LEU:HD12	2:C:63:LEU:C	2.31	0.52
4:C:2151:HOH:O	1:E:179:ARG:HD3	2.09	0.51
4:A:2194:HOH:O	2:C:55:ALA:HB3	2.10	0.51
1:E:211:LYS:O	1:E:214:GLU:HG2	2.11	0.51
1:E:227:ARG:NH2	4:E:2055:HOH:O	2.44	0.51
2:F:63:LEU:C	2:F:63:LEU:HD12	2.30	0.51
1:E:56:LYS:HG2	2:F:200:SER:HA	1.93	0.51
1:B:60:MET:HB2	1:B:178:GLY:HA2	1.94	0.50
2:F:217:THR:HG23	2:F:218:HIS:CE1	2.46	0.50
1:E:63:LEU:C	1:E:63:LEU:HD12	2.33	0.49
1:A:66:ASN:HB2	4:B:2026:HOH:O	2.12	0.49
1:A:90:PRO:HD2	4:A:2043:HOH:O	2.11	0.49
4:D:2129:HOH:O	1:E:116:MET:HE3	2.13	0.49
1:B:178:GLY:O	2:C:205:PRO:HG2	2.13	0.49
1:D:112:GLU:OE2	1:E:227:ARG:NH2	2.45	0.48
2:C:217:THR:HG23	2:C:218:HIS:ND1	2.29	0.48
1:D:12:SER:HB3	1:D:18:PRO:HD3	1.95	0.48
1:A:179:ARG:NH1	1:D:76:ARG:O	2.45	0.48
1:E:178:GLY:O	2:F:205:PRO:HG2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:217:THR:HG22	4:F:349:HOH:O	2.13	0.48
2:C:12:SER:O	2:C:13:GLN:HB2	2.14	0.47
2:C:138:TRP:CZ2	2:C:224:VAL:HG21	2.49	0.47
3:A:2009:ACT:H1	4:B:2170:HOH:O	2.15	0.47
1:B:216:ARG:HD2	4:B:2155:HOH:O	2.13	0.47
1:A:141:TYR:CE2	1:A:216:ARG:HD2	2.49	0.47
1:B:3:ASP:HB3	4:B:2099:HOH:O	2.15	0.47
3:D:2008:ACT:H2	1:E:201:MET:CE	2.45	0.46
1:B:56:LYS:HG2	2:C:200:SER:HA	1.96	0.46
1:D:202:PHE:O	2:F:98:GLU:HB3	2.15	0.46
1:D:20:CYS:SG	1:D:26:ILE:HA	2.56	0.46
2:C:211:LYS:HE3	2:F:93:MET:CG	2.45	0.46
1:B:63:LEU:C	1:B:63:LEU:HD12	2.36	0.46
1:E:72:ASN:ND2	1:E:75:SER:HB2	2.30	0.46
1:A:4:HIS:CE1	1:B:4:HIS:HD2	2.34	0.46
1:A:202:PHE:O	2:C:98:GLU:HB3	2.16	0.46
1:B:160:SER:HB2	4:B:2187:HOH:O	2.15	0.45
2:F:6:TYR:HD2	2:F:113:ALA:HB3	1.81	0.45
1:A:55:ARG:HH12	1:B:128:PRO:HB2	1.82	0.45
1:A:76:ARG:O	1:D:179:ARG:NH1	2.48	0.45
1:B:85:THR:HB	1:B:86:PRO:CD	2.46	0.45
1:D:66:ASN:HB2	4:E:2018:HOH:O	2.16	0.45
1:D:90:PRO:HD2	4:D:2029:HOH:O	2.16	0.45
2:F:12:SER:O	2:F:13:GLN:HB2	2.16	0.45
1:D:131:PRO:HB2	1:D:134:TRP:CD1	2.52	0.45
2:C:6:TYR:HD2	2:C:113:ALA:HB3	1.82	0.45
1:D:99:GLU:HG2	4:E:2171:HOH:O	2.17	0.45
2:F:57:PHE:CD2	2:F:58:SER:N	2.85	0.45
1:A:20:CYS:SG	1:A:26:ILE:HA	2.56	0.44
1:A:145:MET:HG2	1:A:191:PHE:HB2	1.99	0.44
1:B:40:GLU:HG2	2:F:40:GLU:OE2	2.18	0.44
1:D:60:MET:HB2	1:D:178:GLY:HA2	1.99	0.44
1:A:60:MET:HB2	1:A:178:GLY:HA2	1.99	0.44
1:B:141:TYR:OH	1:B:216:ARG:HD3	2.17	0.44
4:A:2117:HOH:O	1:B:116:MET:HE2	2.17	0.44
1:B:40:GLU:OE1	2:F:37:GLU:OE2	2.35	0.44
1:D:55:ARG:HH12	1:E:128:PRO:HB2	1.83	0.44
1:E:60:MET:HB2	1:E:178:GLY:HA2	1.98	0.44
1:E:85:THR:HB	1:E:86:PRO:CD	2.48	0.43
2:F:138:TRP:CE2	2:F:224:VAL:HG21	2.54	0.43
1:B:211:LYS:CE	1:D:93:MET:SD	3.06	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:TYR:OH	1:E:216:ARG:HD3	2.17	0.43
2:F:41:LYS:HE2	4:F:260:HOH:O	2.18	0.43
2:F:131:PRO:HG2	2:F:134:TRP:CG	2.53	0.43
2:F:138:TRP:CZ2	2:F:224:VAL:HG21	2.53	0.43
1:B:175:GLU:O	1:B:182:CYS:HA	2.19	0.43
1:A:40:GLU:OE1	1:D:40:GLU:OE1	2.37	0.43
1:A:138:TRP:CE2	1:A:224:VAL:HG21	2.54	0.43
1:A:12:SER:O	1:A:13:GLN:HB2	2.19	0.43
2:C:40:GLU:OE2	1:E:40:GLU:HG2	2.19	0.43
1:E:175:GLU:O	1:E:182:CYS:HA	2.19	0.43
1:A:71:CYS:SG	1:B:215:LEU:HD22	2.59	0.42
2:C:57:PHE:CD2	2:C:58:SER:N	2.87	0.42
1:D:188:ALA:O	2:F:67:PRO:HD3	2.18	0.42
1:A:122:SER:O	1:A:123:GLN:HB2	2.19	0.42
1:B:66:ASN:OD1	1:B:66:ASN:C	2.57	0.42
2:F:19:MET:CE	2:F:20:CYS:H	2.33	0.42
1:D:145:MET:HG2	1:D:191:PHE:HB2	2.00	0.42
1:A:85:THR:HB	1:A:86:PRO:CD	2.50	0.42
2:F:19:MET:HE3	2:F:20:CYS:H	1.84	0.42
2:F:57:PHE:CG	2:F:58:SER:N	2.88	0.42
1:A:131:PRO:HB2	1:A:134:TRP:CD1	2.55	0.42
2:C:102:LYS:O	2:C:107:ARG:NH2	2.53	0.42
2:C:145:MET:HA	2:C:155:GLY:HA2	2.00	0.42
1:A:55:ARG:NH2	4:A:2189:HOH:O	2.53	0.42
2:F:197:PRO:HG2	2:F:198:GLU:OE2	2.20	0.42
1:E:42:ALA:HB2	1:E:152:GLU:HB3	2.02	0.41
2:F:145:MET:HA	2:F:155:GLY:HA2	2.02	0.41
2:C:131:PRO:HG2	2:C:134:TRP:CG	2.55	0.41
1:E:28:TYR:CZ	1:E:110:VAL:HG21	2.56	0.41
2:C:138:TRP:CE2	2:C:224:VAL:HG21	2.55	0.41
2:F:36:PHE:CE1	2:F:61:PRO:HG2	2.56	0.41
2:C:197:PRO:HG2	2:C:198:GLU:OE2	2.20	0.41
2:F:97:ALA:HB2	2:F:181:THR:HG22	2.02	0.41
1:B:179:ARG:CZ	2:F:178:GLY:HA3	2.51	0.41
2:F:211:LYS:O	2:F:212:ALA:C	2.59	0.41
1:B:17:ASP:HB3	1:B:29:HIS:CD2	2.56	0.41
2:C:122:SER:O	2:C:123:GLN:HB2	2.21	0.41
1:D:12:SER:O	1:D:13:GLN:HB2	2.20	0.41
1:E:17:ASP:HB3	1:E:29:HIS:CD2	2.56	0.41
4:E:2156:HOH:O	2:F:227:LYS:NZ	2.54	0.41
1:A:93:MET:SD	1:E:211:LYS:CE	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ARG:HH11	1:A:41:ARG:HG2	1.86	0.40
1:D:85:THR:HB	1:D:86:PRO:CD	2.51	0.40
1:A:188:ALA:O	2:C:67:PRO:HD3	2.21	0.40
3:A:2009:ACT:H2	1:B:201:MET:CE	2.52	0.40
1:B:12:SER:O	1:B:13:GLN:HB2	2.20	0.40
2:C:107:ARG:HA	2:C:107:ARG:HD3	1.91	0.40
2:C:211:LYS:O	2:C:212:ALA:C	2.59	0.40
2:F:125:VAL:O	2:F:221:ARG:NE	2.53	0.40
1:B:198:ARG:NH1	4:B:2170:HOH:O	2.55	0.40
2:C:17:GLU:OE1	2:C:26:LYS:NZ	2.47	0.40
1:D:141:TYR:CE2	1:D:216:ARG:HD2	2.57	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	226/229 (99%)	221 (98%)	3 (1%)	2 (1%)	17	7
1	B	225/229 (98%)	217 (96%)	8 (4%)	0	100	100
1	D	226/229 (99%)	221 (98%)	3 (1%)	2 (1%)	17	7
1	E	225/229 (98%)	218 (97%)	7 (3%)	0	100	100
2	C	222/228 (97%)	210 (95%)	12 (5%)	0	100	100
2	F	223/228 (98%)	211 (95%)	12 (5%)	0	100	100
All	All	1347/1372 (98%)	1298 (96%)	45 (3%)	4 (0%)	41	31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	SER

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Mol	Chain	Res	Type
1	A	76	ARG
1	D	76	ARG
1	D	148	SER

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	194/195 (100%)	192 (99%)	2 (1%)	76	76
1	B	193/195 (99%)	190 (98%)	3 (2%)	62	60
1	D	194/195 (100%)	192 (99%)	2 (1%)	76	76
1	E	193/195 (99%)	192 (100%)	1 (0%)	88	89
2	C	188/192 (98%)	186 (99%)	2 (1%)	73	73
2	F	189/192 (98%)	186 (98%)	3 (2%)	62	60
All	All	1151/1164 (99%)	1138 (99%)	13 (1%)	73	73

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	MET
1	A	227	ARG
1	B	72	ASN
1	B	89	MET
1	B	227	ARG
2	C	66	ASN
2	C	99	ASP
1	D	145	MET
1	D	227	ARG
1	E	89	MET
2	F	66	ASN
2	F	99	ASP
2	F	169	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	29	HIS
1	B	4	HIS
1	B	29	HIS
2	C	25	ASN
2	C	202	GLN
1	D	19	GLN
1	D	29	HIS
1	E	4	HIS
1	E	29	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 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	ACT	D	2008	-	3,3,3	0.94	0	3,3,3	0.52	0
3	ACT	B	2007	-	3,3,3	0.85	0	3,3,3	0.88	0
3	ACT	C	2004	-	3,3,3	0.98	0	3,3,3	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACT	B	2002	-	3,3,3	0.92	0	3,3,3	0.77	0
3	ACT	B	2006	-	3,3,3	0.94	0	3,3,3	0.66	0
3	ACT	B	2003	-	3,3,3	0.92	0	3,3,3	0.70	0
3	ACT	A	2009	-	3,3,3	1.00	0	3,3,3	0.49	0
3	ACT	E	2001	-	3,3,3	0.89	0	3,3,3	0.74	0
3	ACT	C	2005	-	3,3,3	0.91	0	3,3,3	0.72	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2008	ACT	1	0
3	A	2009	ACT	2	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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.