



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 14, 2017 – 10:26 PM EDT

PDB ID : 1SHL
Title : CASPASE-7 IN COMPLEX WITH FICA ALLOSTERIC INHIBITOR
Authors : Hardy, J.A.; Lam, J.; Nguyen, J.T.; O'Brien, T.; Wells, J.A.
Deposited on : unknown
Resolution : 3.00 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

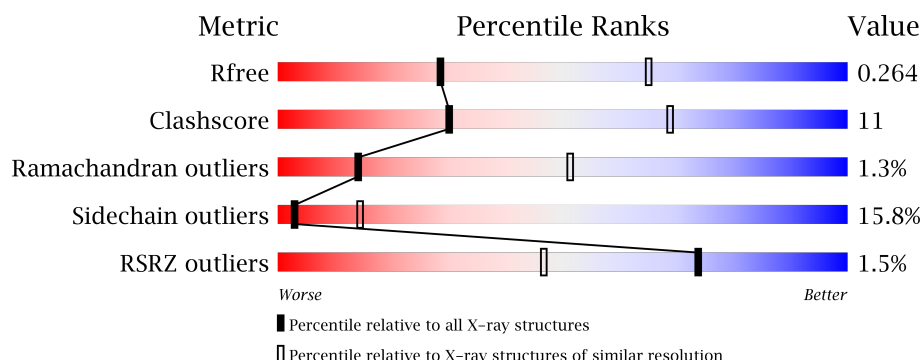
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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 ≥ 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.

Mol	Chain	Length	Quality of chain
1	A	245	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 53%, yellow 22%, orange 5%, grey 19%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 100% 53% 22% • • 19% </div> </div>
1	B	245	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 53%, yellow 23%, orange 5%, grey 18%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 100% 53% 23% 5% 18% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3196 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.

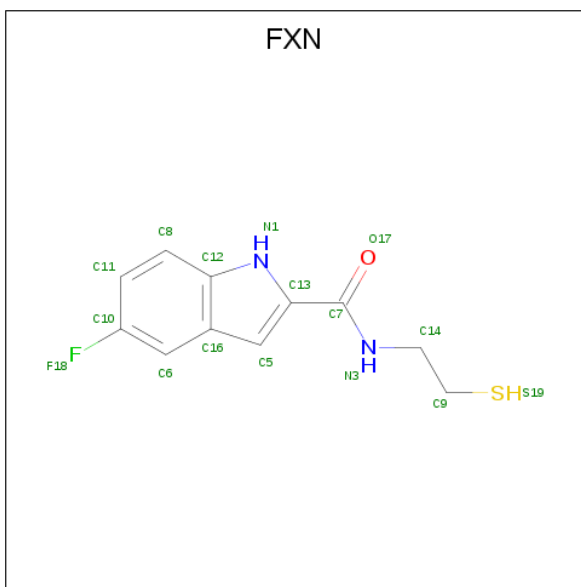
- Molecule 1 is a protein called Caspase-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	8	0	0
			1576	1006	266	289	15			
1	B	200	Total	C	N	O	S	4	0	0
			1588	1013	268	292	15			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	192	ALA	ASP	ENGINEERED	UNP P55210
A	304	GLN	-	EXPRESSION TAG	UNP P55210
A	305	LEU	-	EXPRESSION TAG	UNP P55210
A	306	HIS	-	EXPRESSION TAG	UNP P55210
A	307	HIS	-	EXPRESSION TAG	UNP P55210
A	308	HIS	-	EXPRESSION TAG	UNP P55210
A	309	HIS	-	EXPRESSION TAG	UNP P55210
A	310	HIS	-	EXPRESSION TAG	UNP P55210
A	311	HIS	-	EXPRESSION TAG	UNP P55210
B	192	ALA	ASP	ENGINEERED	UNP P55210
B	304	GLN	-	EXPRESSION TAG	UNP P55210
B	305	LEU	-	EXPRESSION TAG	UNP P55210
B	306	HIS	-	EXPRESSION TAG	UNP P55210
B	307	HIS	-	EXPRESSION TAG	UNP P55210
B	308	HIS	-	EXPRESSION TAG	UNP P55210
B	309	HIS	-	EXPRESSION TAG	UNP P55210
B	310	HIS	-	EXPRESSION TAG	UNP P55210
B	311	HIS	-	EXPRESSION TAG	UNP P55210

- Molecule 2 is 5-FLUORO-1H-INDOLE-2-CARBOXYLIC ACID-(2-MERCAPTO-ETHYL)-AMIDE (three-letter code: FXN) (formula: C₁₁H₁₁FN₂OS).

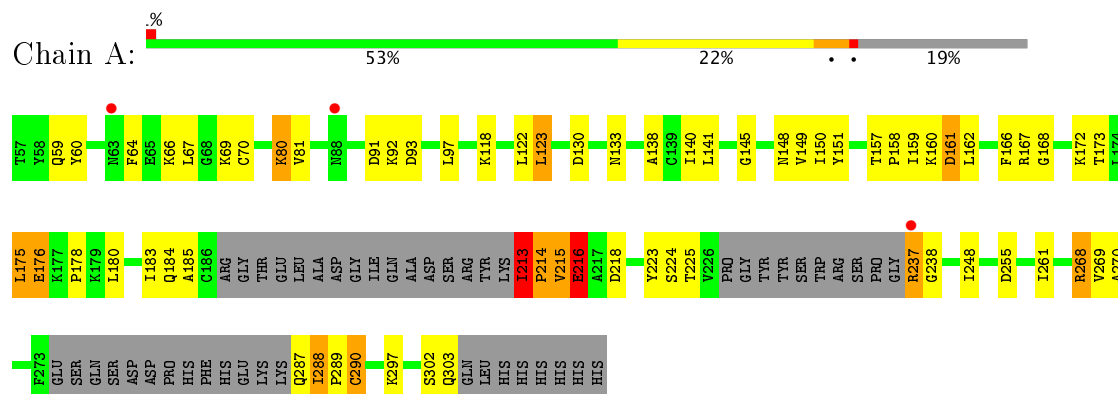


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	F	N	O	S		
2	A	1	16	11	1	2	1	1	0	0
2	B	1	16	11	1	2	1	1	0	0

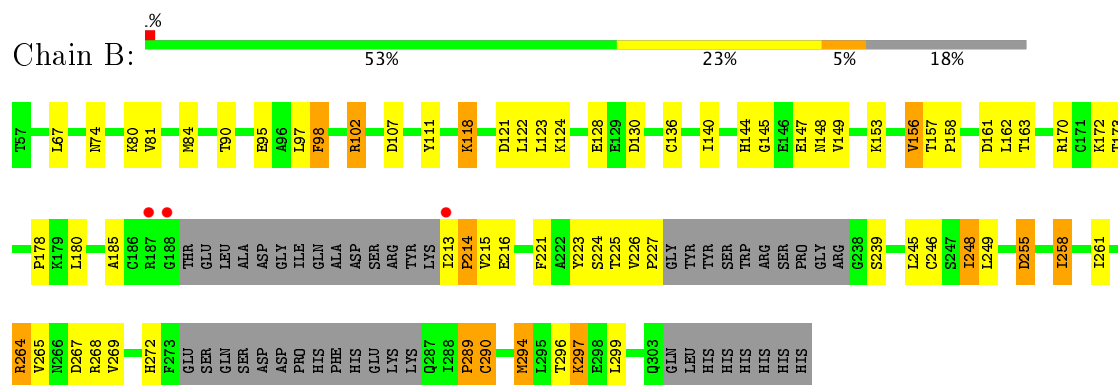
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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-7



• Molecule 1: Caspase-7



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	90.22Å 90.22Å 186.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 3.00 19.97 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (10.00-3.00) 99.1 (19.97-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.60 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.221 , 0.273 0.214 , 0.264	Depositor DCC
R_{free} test set	897 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	60.8	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3196	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: FXN

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	0.52	2/1603 (0.1%)	0.82	8/2152 (0.4%)
1	B	0.42	0/1616	0.79	6/2171 (0.3%)
All	All	0.47	2/3219 (0.1%)	0.80	14/4323 (0.3%)

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 (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	LYS	CB-CG	-8.99	1.28	1.52
1	A	290	CYS	C-N	-5.74	1.20	1.34

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	GLN	CA-CB-CG	-10.60	90.09	113.40
1	A	130	ASP	CB-CG-OD2	6.35	124.01	118.30
1	A	290	CYS	O-C-N	-6.25	112.71	122.70
1	A	172	LYS	CB-CG-CD	-5.97	96.08	111.60
1	B	130	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	218	ASP	CB-CG-OD2	5.55	123.29	118.30
1	B	267	ASP	CB-CG-OD2	5.36	123.13	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	ASP	CB-CG-OD2	5.34	123.10	118.30
1	A	91	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	289	PRO	O-C-N	-5.27	114.27	122.70
1	A	255	ASP	CB-CG-OD2	5.25	123.03	118.30
1	A	161	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	121	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	255	ASP	CB-CG-OD2	5.09	122.88	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	213	ILE	Peptide
1	B	290	CYS	Mainchain

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	1576	0	1565	32	0
1	B	1588	0	1576	35	0
2	A	16	0	10	1	0
2	B	16	0	10	1	0
All	All	3196	0	3161	67	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 (67) 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:145:GLY:HA3	1:A:185:ALA:HB1	1.55	0.89
1:B:111:TYR:CZ	1:B:122:LEU:HD11	2.12	0.85
1:A:223:TYR:CD1	2:B:401:FXN:H11	2.11	0.85
1:A:213:ILE:HB	1:A:214:PRO:HA	1.60	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:VAL:O	1:A:216:GLU:HB2	1.83	0.77
1:B:214:PRO:HG2	1:B:221:PHE:CZ	2.24	0.73
1:B:124:LYS:HE3	1:B:128:GLU:OE2	1.92	0.70
1:B:111:TYR:CE2	1:B:122:LEU:HD11	2.27	0.68
1:B:249:LEU:HD13	1:B:261:ILE:HG21	1.75	0.68
1:A:141:LEU:HD12	1:A:183:ILE:HG12	1.81	0.63
2:A:401:FXN:H11	1:B:223:TYR:CD2	2.33	0.63
1:A:175:LEU:HD23	1:A:176:GLU:HG2	1.81	0.63
1:A:213:ILE:CB	1:A:214:PRO:HA	2.29	0.61
1:B:118:LYS:O	1:B:122:LEU:HB2	2.02	0.60
1:B:97:LEU:HD13	1:B:140:ILE:HG21	1.83	0.59
1:A:93:ASP:OD1	1:A:238:GLY:HA3	2.02	0.58
1:A:158:PRO:HG2	1:A:161:ASP:OD2	2.05	0.57
1:A:223:TYR:HB3	1:A:290:CYS:SG	2.45	0.57
1:A:92:LYS:HD3	1:A:237:ARG:HA	1.86	0.57
1:A:123:LEU:HD23	1:A:166:PHE:HE2	1.70	0.56
1:A:248:ILE:HD11	1:A:268:ARG:NH1	2.20	0.56
1:A:123:LEU:HD23	1:A:166:PHE:CE2	2.42	0.55
1:A:70:CYS:HA	1:A:138:ALA:O	2.07	0.55
1:B:213:ILE:HG23	1:B:215:VAL:HG23	1.87	0.55
1:A:148:ASN:H	1:A:149:VAL:HG12	1.72	0.52
1:B:248:ILE:HD12	1:B:264:ARG:HB3	1.91	0.52
1:B:269:VAL:CG2	1:B:289:PRO:HD3	2.40	0.51
1:B:269:VAL:HG21	1:B:289:PRO:HD3	1.93	0.51
1:B:215:VAL:HA	1:B:294:MET:HE2	1.93	0.50
1:B:74:ASN:HD21	1:B:90:THR:HG23	1.76	0.50
1:A:150:ILE:HG22	1:A:151:TYR:N	2.26	0.50
1:B:215:VAL:HA	1:B:294:MET:CE	2.41	0.50
1:A:268:ARG:HG2	1:A:268:ARG:HH11	1.77	0.49
1:B:123:LEU:HD12	1:B:162:LEU:HD22	1.94	0.49
1:B:213:ILE:CG2	1:B:215:VAL:HG23	2.43	0.48
1:A:268:ARG:HG2	1:A:268:ARG:NH1	2.29	0.48
1:A:159:ILE:HD12	1:A:183:ILE:HD13	1.96	0.48
1:B:163:THR:HG21	1:B:221:PHE:HE1	1.79	0.47
1:A:215:VAL:O	1:A:216:GLU:CB	2.60	0.47
1:A:97:LEU:HD13	1:A:140:ILE:HG21	1.96	0.47
1:A:270:ALA:HB2	1:A:288:ILE:HD12	1.96	0.47
1:B:98:PHE:O	1:B:102:ARG:HB2	2.15	0.47
1:A:148:ASN:OD1	1:A:148:ASN:N	2.47	0.47
1:B:149:VAL:CG1	1:B:156:VAL:HG13	2.45	0.46
1:A:64:PHE:HB3	1:A:133:ASN:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:LEU:HD23	1:B:246:CYS:SG	2.56	0.46
1:B:223:TYR:HB3	1:B:290:CYS:SG	2.57	0.45
1:B:258:ILE:CG1	1:B:299:LEU:HB3	2.47	0.45
1:B:84:MET:HB3	1:B:144:HIS:CD2	2.51	0.45
1:B:258:ILE:H	1:B:258:ILE:HG13	1.60	0.44
1:A:167:ARG:HD3	1:A:216:GLU:OE1	2.17	0.44
1:B:245:LEU:O	1:B:249:LEU:HB2	2.18	0.43
1:B:261:ILE:O	1:B:265:VAL:HG23	2.19	0.43
1:B:122:LEU:HD12	1:B:122:LEU:HA	1.81	0.43
1:B:296:THR:C	1:B:297:LYS:HG2	2.39	0.43
1:A:168:GLY:O	1:A:175:LEU:HD12	2.19	0.42
1:B:136:CYS:HB3	1:B:178:PRO:HG2	2.02	0.42
1:A:268:ARG:CG	1:A:268:ARG:HH11	2.32	0.42
1:A:123:LEU:HD13	1:A:162:LEU:HD22	2.01	0.42
1:B:145:GLY:HA3	1:B:185:ALA:HB1	2.01	0.42
1:A:184:GLN:O	1:A:184:GLN:HG3	2.19	0.41
1:B:213:ILE:HG23	1:B:215:VAL:CG2	2.50	0.41
1:A:60:TYR:CD1	1:A:178:PRO:HD3	2.55	0.41
1:B:158:PRO:HG2	1:B:161:ASP:OD2	2.21	0.40
1:A:288:ILE:HG13	1:A:289:PRO:HD2	2.03	0.40
1:B:149:VAL:HG11	1:B:156:VAL:HG13	2.02	0.40
1:B:226:VAL:HA	1:B:227:PRO:HD3	1.89	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	190/245 (78%)	175 (92%)	11 (6%)	4 (2%)	8	38
1	B	192/245 (78%)	177 (92%)	14 (7%)	1 (0%)	32	74
All	All	382/490 (78%)	352 (92%)	25 (6%)	5 (1%)	14	51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	LYS
1	A	216	GLU
1	A	214	PRO
1	B	214	PRO
1	A	81	VAL

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	174/215 (81%)	147 (84%)	27 (16%)	3	15
1	B	175/215 (81%)	147 (84%)	28 (16%)	3	14
All	All	349/430 (81%)	294 (84%)	55 (16%)	3	14

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	66	LYS
1	A	67	LEU
1	A	69	LYS
1	A	80	LYS
1	A	118	LYS
1	A	122	LEU
1	A	123	LEU
1	A	157	THR
1	A	160	LYS
1	A	173	THR
1	A	175	LEU
1	A	176	GLU
1	A	180	LEU
1	A	213	ILE
1	A	215	VAL
1	A	216	GLU
1	A	224	SER

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Mol	Chain	Res	Type
1	A	225	THR
1	A	237	ARG
1	A	261	ILE
1	A	268	ARG
1	A	269	VAL
1	A	288	ILE
1	A	297	LYS
1	A	302	SER
1	A	303	GLN
1	B	67	LEU
1	B	80	LYS
1	B	81	VAL
1	B	95	GLU
1	B	98	PHE
1	B	102	ARG
1	B	118	LYS
1	B	147	GLU
1	B	148	ASN
1	B	153	LYS
1	B	156	VAL
1	B	157	THR
1	B	170	ARG
1	B	172	LYS
1	B	173	THR
1	B	180	LEU
1	B	216	GLU
1	B	224	SER
1	B	225	THR
1	B	239	SER
1	B	248	ILE
1	B	255	ASP
1	B	258	ILE
1	B	264	ARG
1	B	268	ARG
1	B	272	HIS
1	B	294	MET
1	B	297	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN

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Mol	Chain	Res	Type
1	A	112	ASN
1	A	144	HIS
1	B	74	ASN
1	B	112	ASN
1	B	148	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](#)

2 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	$\# Z > 2$
2	FXN	A	401	1	15,17,17	1.20	2 (13%)	17,23,23	1.74	3 (17%)
2	FXN	B	401	1	15,17,17	1.19	1 (6%)	17,23,23	1.86	3 (17%)

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
2	FXN	A	401	1	-	0/5/8/8	0/2/2/2
2	FXN	B	401	1	-	0/5/8/8	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FXN	C11-C10	2.05	1.41	1.37
2	A	401	FXN	C6-C10	2.91	1.40	1.36
2	B	401	FXN	C6-C10	3.31	1.41	1.36

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FXN	C11-C8-C12	-3.63	116.39	120.84
2	B	401	FXN	C11-C8-C12	-3.15	116.97	120.84
2	A	401	FXN	C8-C11-C10	2.24	121.34	118.70
2	B	401	FXN	C8-C11-C10	2.48	121.62	118.70
2	A	401	FXN	C13-C7-N3	3.65	120.52	115.52
2	B	401	FXN	C13-C7-N3	4.60	121.83	115.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FXN	1	0
2	B	401	FXN	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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 ²)	Q<0.9
1	A	198/245 (80%)	-0.40	3 (1%)	74 47	35, 55, 85, 107	2 (1%)
1	B	200/245 (81%)	-0.42	3 (1%)	74 47	35, 54, 83, 100	1 (0%)
All	All	398/490 (81%)	-0.41	6 (1%)	74 47	35, 54, 85, 107	3 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	188	GLY	5.1
1	A	88	ASN	2.3
1	B	213	ILE	2.3
1	A	237	ARG	2.3
1	A	63	ASN	2.2
1	B	187	ARG	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	LLDF	B-factors(\AA^2)	Q<0.9
2	FXN	B	401	16/16	0.93	0.25	0.95	43,46,54,61	0
2	FXN	A	401	16/16	0.89	0.30	0.66	60,62,66,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.