



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:54 AM UTC

PDB ID : 6OJA / pdb_00006oja
Title : Crystal structure of the N. meningitides methionine-binding protein in its L-methionine bound conformation
Authors : Nguyen, P.T.; Lai, J.Y.; Kaiser, J.T.; Rees, D.C.
Deposited on : 2019-04-11
Resolution : 1.55 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

i

X-RAY DIFFRACTION

A.

Metric	Percentile Ranks	Value
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Percentile relative to all X-ray structures

Percentile relative to X-ray structures of similar resolution

Metric

Whole archive
(#Entries)

Similar resolution
(#Entries, resolution range(Å))

 R_{free}

Clashscore

Ramachandran outliers

Sidechain outliers

RSRZ outliers

180053

190562

187476

187428

180081

2145 (1.56-1.56)

2189 (1.56-1.56)

2153 (1.56-1.56)

2150 (1.56-1.56)

2146 (1.56-1.56)

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 $\leq 5\%$. 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.

Quality of chain

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Mol	Chain	Length	Quality of chain
1	F	248	3% 92% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 24411 atoms, of which 11238 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 Lipoprotein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			
1	B	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			
1	C	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			
1	D	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			
1	E	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			
1	F	241	Total	C	H	N	O	S	0	0	0
			3764	1218	1865	308	370	3			

There are 18 discrepancies between the modelled and reference sequences:

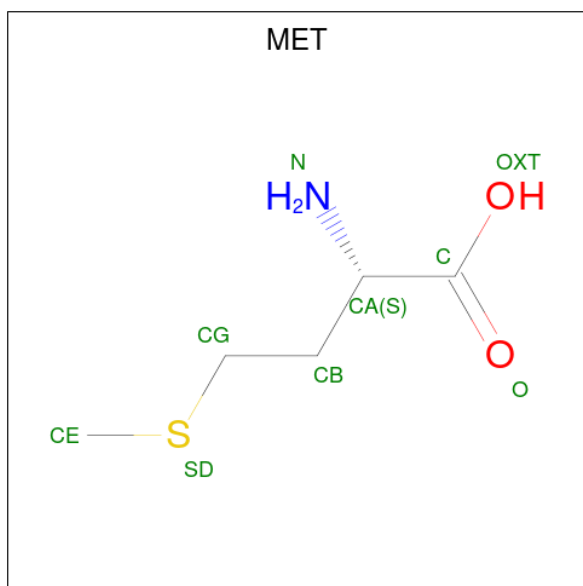
Chain	Residue	Modelled	Actual	Comment	Reference
A	40	LYS	-	expression tag	UNP Q9JPG4
A	41	HIS	-	expression tag	UNP Q9JPG4
A	42	MET	-	expression tag	UNP Q9JPG4
B	40	LYS	-	expression tag	UNP Q9JPG4
B	41	HIS	-	expression tag	UNP Q9JPG4
B	42	MET	-	expression tag	UNP Q9JPG4
C	40	LYS	-	expression tag	UNP Q9JPG4
C	41	HIS	-	expression tag	UNP Q9JPG4
C	42	MET	-	expression tag	UNP Q9JPG4
D	40	LYS	-	expression tag	UNP Q9JPG4
D	41	HIS	-	expression tag	UNP Q9JPG4
D	42	MET	-	expression tag	UNP Q9JPG4
E	40	LYS	-	expression tag	UNP Q9JPG4
E	41	HIS	-	expression tag	UNP Q9JPG4
E	42	MET	-	expression tag	UNP Q9JPG4
F	40	LYS	-	expression tag	UNP Q9JPG4
F	41	HIS	-	expression tag	UNP Q9JPG4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	42	MET	-	expression tag	UNP Q9JPG4

- Molecule 2 is METHIONINE (CCD ID: MET) (formula: $C_5H_{11}NO_2S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		
2	B	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		
2	C	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		
2	D	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		
2	E	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		
2	F	1	Total	C	H	N	O	S	0	0
			17	5	8	1	2	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	301	Total	O	0	0
			301	301		
3	B	288	Total	O	0	0
			288	288		
3	C	258	Total	O	0	0
			258	258		

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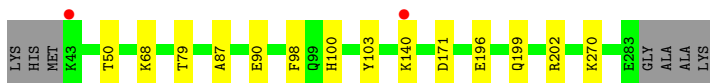
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	304	Total 304	O 304	0	0
3	E	281	Total 281	O 281	0	0
3	F	293	Total 293	O 293	0	0

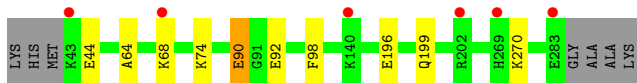
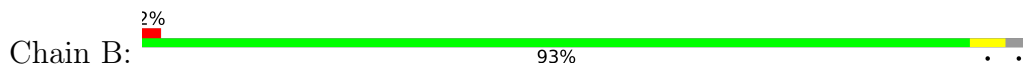
3 Residue-property plots [i](#)

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 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: Lipoprotein



- Molecule 1: Lipoprotein



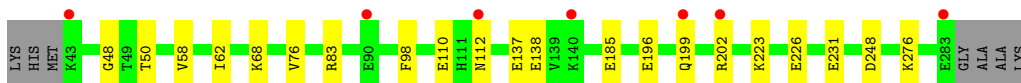
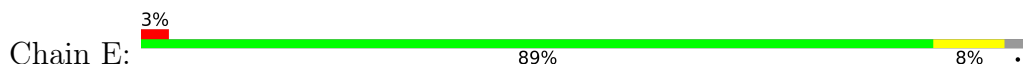
- Molecule 1: Lipoprotein



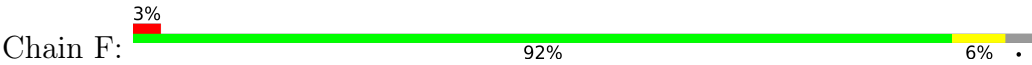
- Molecule 1: Lipoprotein



- Molecule 1: Lipoprotein



- Molecule 1: Lipoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.58Å 87.65Å 91.63Å 114.70° 104.41° 105.24°	Depositor
Resolution (Å)	34.20 – 1.55 34.20 – 1.55	Depositor EDS
% Data completeness (in resolution range)	96.0 (34.20-1.55) 96.0 (34.20-1.55)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.55Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.162 , 0.181 0.166 , 0.183	Depositor DCC
R_{free} test set	13993 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	24411	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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 [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	0.37	0/1941	0.56	0/2633
1	B	0.36	0/1941	0.54	0/2633
1	C	0.37	0/1941	0.54	0/2633
1	D	0.36	0/1941	0.56	0/2633
1	E	0.37	0/1941	0.55	0/2633
1	F	0.38	0/1941	0.56	0/2633
All	All	0.37	0/11646	0.55	0/15798

There are no bond length outliers.

There are no bond angle outliers.

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	1899	1865	1865	11	0
1	B	1899	1865	1865	8	0
1	C	1899	1865	1865	20	0
1	D	1899	1865	1865	12	0
1	E	1899	1865	1865	19	0
1	F	1899	1865	1865	18	0
2	A	9	8	8	0	0
2	B	9	8	8	0	0
2	C	9	8	8	0	0
2	D	9	8	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	9	8	8	0	0
2	F	9	8	8	0	0
3	A	301	0	0	5	4
3	B	288	0	0	5	0
3	C	258	0	0	11	3
3	D	304	0	0	9	0
3	E	281	0	0	9	3
3	F	293	0	0	11	4
All	All	13173	11238	11238	75	7

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 3.

All (75) 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:F:202:ARG:NH2	3:F:1101:HOH:O	1.86	1.07
1:D:112:ASN:CG	3:D:1101:HOH:O	2.19	0.83
1:E:112:ASN:ND2	1:F:199:GLN:OE1	2.12	0.82
1:F:90:GLU:OE1	3:F:1102:HOH:O	1.98	0.81
1:E:138:GLU:OE2	3:E:1101:HOH:O	1.98	0.81
1:D:112:ASN:OD1	3:D:1101:HOH:O	2.00	0.80
1:D:137:GLU:OE2	3:D:1102:HOH:O	2.00	0.79
1:F:83:ARG:NH2	3:F:1104:HOH:O	2.16	0.79
1:E:248:ASP:OD1	3:E:1102:HOH:O	2.03	0.76
1:C:68:LYS:NZ	3:C:1102:HOH:O	2.17	0.76
1:C:110:GLU:OE1	3:C:1101:HOH:O	2.03	0.75
1:D:112:ASN:HB3	3:D:1101:HOH:O	1.88	0.73
1:B:196:GLU:OE1	3:B:1101:HOH:O	2.09	0.70
1:A:171:ASP:OD1	3:A:1101:HOH:O	2.08	0.69
1:B:44:GLU:OE1	3:B:1102:HOH:O	2.10	0.69
1:C:112:ASN:ND2	3:C:1103:HOH:O	2.25	0.68
1:C:199:GLN:NE2	3:D:1101:HOH:O	2.27	0.67
1:D:112:ASN:CB	3:D:1101:HOH:O	2.40	0.67
1:E:202:ARG:NH2	3:E:1105:HOH:O	2.27	0.67
1:D:90:GLU:OE1	3:D:1103:HOH:O	2.12	0.66
1:B:199:GLN:NE2	3:B:1103:HOH:O	2.15	0.66
1:E:137:GLU:HG3	3:E:1106:HOH:O	1.94	0.66
1:F:110:GLU:OE1	3:F:1103:HOH:O	2.14	0.65
1:C:202:ARG:CZ	3:C:1106:HOH:O	2.46	0.64
1:E:226:GLU:OE2	3:E:1103:HOH:O	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:ASN:HD22	1:D:199:GLN:CD	2.06	0.63
1:C:202:ARG:NH1	3:C:1106:HOH:O	2.32	0.62
1:E:68:LYS:HG2	3:E:1220:HOH:O	1.98	0.62
1:C:282:ASN:ND2	3:C:1105:HOH:O	2.32	0.62
1:A:140:LYS:HE2	3:A:1129:HOH:O	1.98	0.62
1:C:112:ASN:N	1:C:112:ASN:OD1	2.36	0.58
1:E:223:LYS:HB2	1:E:226:GLU:HG3	1.85	0.58
1:F:168:LYS:HE3	3:F:1215:HOH:O	2.03	0.58
1:E:231:GLU:O	1:E:276:LYS:NZ	2.38	0.57
1:A:68:LYS:NZ	3:A:1108:HOH:O	2.38	0.56
1:E:112:ASN:ND2	1:F:199:GLN:CD	2.64	0.55
1:E:68:LYS:HE3	3:E:1220:HOH:O	2.06	0.54
1:F:83:ARG:CD	3:F:1274:HOH:O	2.55	0.54
1:A:270:LYS:NZ	1:C:204:ARG:HH22	2.06	0.54
1:C:270:LYS:CE	3:C:1108:HOH:O	2.56	0.52
1:D:270:LYS:NZ	3:D:1105:HOH:O	2.33	0.52
1:E:112:ASN:ND2	1:F:196:GLU:HB2	2.26	0.51
1:C:112:ASN:CG	1:D:196:GLU:HB2	2.35	0.51
1:F:132:LYS:NZ	3:F:1106:HOH:O	2.44	0.51
1:C:270:LYS:NZ	3:C:1108:HOH:O	2.39	0.51
1:E:199:GLN:CD	1:F:112:ASN:ND2	2.70	0.50
1:A:79:THR:HA	3:A:1191:HOH:O	2.12	0.49
1:D:112:ASN:ND2	3:D:1110:HOH:O	2.46	0.48
1:B:74:LYS:NZ	3:B:1105:HOH:O	2.31	0.47
1:F:83:ARG:HD3	3:F:1274:HOH:O	2.13	0.47
1:C:282:ASN:CG	3:C:1105:HOH:O	2.57	0.47
1:B:64:ALA:O	1:B:68:LYS:HG3	2.15	0.46
1:C:83:ARG:NH2	1:D:92:GLU:OE2	2.49	0.46
1:C:112:ASN:OD1	1:D:196:GLU:HB2	2.16	0.46
1:F:83:ARG:HD2	3:F:1274:HOH:O	2.16	0.46
1:C:270:LYS:HE3	3:C:1108:HOH:O	2.16	0.46
1:F:140:LYS:HE3	3:F:1124:HOH:O	2.17	0.45
1:F:100:HIS:CE1	1:F:103:TYR:HB2	2.52	0.44
1:C:202:ARG:NH2	3:C:1106:HOH:O	2.48	0.44
1:B:196:GLU:CD	3:B:1101:HOH:O	2.59	0.44
1:A:199:GLN:HG2	1:A:202:ARG:HH22	1.83	0.44
1:E:196:GLU:HB2	1:F:112:ASN:ND2	2.34	0.43
1:A:87:ALA:HA	1:B:90:GLU:OE2	2.18	0.43
1:E:83:ARG:NE	3:E:1111:HOH:O	2.51	0.43
1:E:110:GLU:OE1	3:E:1104:HOH:O	2.21	0.42
1:A:270:LYS:HB3	1:A:270:LYS:HE3	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:48:GLY:HA2	1:E:76:VAL:O	2.20	0.42
1:C:148:PRO:HD2	1:C:157:VAL:HG21	2.02	0.41
1:A:90:GLU:OE2	1:B:92:GLU:OE2	2.38	0.41
1:C:58:VAL:HA	1:C:62:ILE:HB	2.02	0.41
1:F:202:ARG:HG3	3:F:1275:HOH:O	2.20	0.41
1:E:112:ASN:HD22	1:F:199:GLN:CD	2.29	0.41
1:A:100:HIS:CE1	1:A:103:TYR:HB2	2.56	0.40
1:E:58:VAL:HA	1:E:62:ILE:HB	2.03	0.40
1:A:196:GLU:CD	3:A:1127:HOH:O	2.64	0.40

All (7) 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)
3:A:1227:HOH:O	3:F:1292:HOH:O[1_545]	1.75	0.45
3:C:1218:HOH:O	3:E:1231:HOH:O[1_556]	1.89	0.31
3:A:1392:HOH:O	3:F:1345:HOH:O[1_545]	1.98	0.22
3:C:1295:HOH:O	3:E:1176:HOH:O[1_556]	2.07	0.13
3:A:1332:HOH:O	3:F:1300:HOH:O[1_545]	2.10	0.10
3:A:1330:HOH:O	3:F:1235:HOH:O[1_545]	2.12	0.08
3:C:1319:HOH:O	3:E:1179:HOH:O[1_556]	2.13	0.07

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	239/248 (96%)	233 (98%)	5 (2%)	1 (0%)	30 13
1	B	239/248 (96%)	234 (98%)	4 (2%)	1 (0%)	30 13
1	C	239/248 (96%)	233 (98%)	5 (2%)	1 (0%)	30 13
1	D	239/248 (96%)	233 (98%)	5 (2%)	1 (0%)	30 13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	239/248 (96%)	233 (98%)	5 (2%)	1 (0%)	30	13
1	F	239/248 (96%)	233 (98%)	5 (2%)	1 (0%)	30	13
All	All	1434/1488 (96%)	1399 (98%)	29 (2%)	6 (0%)	30	13

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	98	PHE
1	A	98	PHE
1	C	98	PHE
1	D	98	PHE
1	E	98	PHE
1	F	98	PHE

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	203/208 (98%)	202 (100%)	1 (0%)	81	68
1	B	203/208 (98%)	201 (99%)	2 (1%)	68	47
1	C	203/208 (98%)	202 (100%)	1 (0%)	81	68
1	D	203/208 (98%)	202 (100%)	1 (0%)	81	68
1	E	203/208 (98%)	201 (99%)	2 (1%)	68	47
1	F	203/208 (98%)	202 (100%)	1 (0%)	81	68
All	All	1218/1248 (98%)	1210 (99%)	8 (1%)	76	59

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	THR
1	B	90	GLU
1	B	270	LYS
1	C	112	ASN

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Mol	Chain	Res	Type
1	D	137	GLU
1	E	50	THR
1	E	185	GLU
1	F	137	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	213	ASN
1	A	238	ASN
1	B	213	ASN
1	B	238	ASN
1	C	61	GLN
1	C	112	ASN
1	C	199	GLN
1	C	238	ASN
1	D	112	ASN
1	D	199	GLN
1	D	238	ASN
1	D	282	ASN
1	E	112	ASN
1	E	199	GLN
1	E	238	ASN
1	F	112	ASN
1	F	199	GLN
1	F	230	GLN
1	F	238	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 oligosaccharides in this entry.

5.6 Ligand geometry

6 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	MET	D	1001	-	7,8,8	0.79	0	5,9,9	0.45	0
2	MET	A	1001	-	7,8,8	0.71	0	5,9,9	1.03	1 (20%)
2	MET	B	1001	-	7,8,8	0.95	1 (14%)	5,9,9	0.86	0
2	MET	C	1001	-	7,8,8	0.67	0	5,9,9	0.78	0
2	MET	F	1001	-	7,8,8	0.94	1 (14%)	5,9,9	0.87	0
2	MET	E	1001	-	7,8,8	0.88	1 (14%)	5,9,9	0.58	0

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	MET	D	1001	-	-	0/8/8/8	-
2	MET	A	1001	-	-	2/8/8/8	-
2	MET	B	1001	-	-	0/8/8/8	-
2	MET	C	1001	-	-	0/8/8/8	-
2	MET	F	1001	-	-	1/8/8/8	-
2	MET	E	1001	-	-	0/8/8/8	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	MET	OXT-C	-2.27	1.23	1.30
2	F	1001	MET	OXT-C	-2.20	1.23	1.30
2	E	1001	MET	OXT-C	-2.06	1.24	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	1001	MET	OXT-C-O	-2.18	119.13	124.08

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	MET	OXT-C-CA-N
2	F	1001	MET	OXT-C-CA-N
2	A	1001	MET	O-C-CA-N

There are no ring outliers.

No monomer is involved in short contacts.

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

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	241/248 (97%)	-0.04	2 (0%) 82 87	20, 27, 43, 58	0
1	B	241/248 (97%)	0.07	6 (2%) 58 66	20, 29, 44, 62	0
1	C	241/248 (97%)	0.14	9 (3%) 45 53	20, 29, 46, 67	0
1	D	241/248 (97%)	0.08	6 (2%) 58 66	20, 28, 44, 56	0
1	E	241/248 (97%)	0.14	7 (2%) 53 62	21, 29, 45, 68	0
1	F	241/248 (97%)	0.05	8 (3%) 49 57	20, 28, 43, 57	0
All	All	1446/1488 (97%)	0.08	38 (2%) 57 65	20, 28, 45, 68	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	112	ASN	6.0
1	F	112	ASN	5.6
1	C	283	GLU	5.0
1	C	282	ASN	4.5
1	E	112	ASN	3.9
1	D	43	LYS	3.4
1	D	202	ARG	3.3
1	B	68	LYS	3.2
1	B	269	HIS	3.1
1	C	202	ARG	3.0
1	E	283	GLU	2.8
1	E	202	ARG	2.8
1	F	43	LYS	2.8
1	B	43	LYS	2.7
1	A	43	LYS	2.6
1	C	172	GLY	2.6
1	D	199	GLN	2.6
1	C	234	PHE	2.5
1	A	140	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	282	ASN	2.4
1	E	140	LYS	2.4
1	E	43	LYS	2.4
1	B	140	LYS	2.3
1	F	202	ARG	2.2
1	F	72	THR	2.2
1	C	43	LYS	2.2
1	D	112	ASN	2.2
1	F	137	GLU	2.2
1	D	72	THR	2.2
1	B	202	ARG	2.1
1	C	270	LYS	2.1
1	B	283	GLU	2.1
1	E	90	GLU	2.1
1	F	260	ASN	2.1
1	F	283	GLU	2.1
1	C	68	LYS	2.1
1	D	247	LYS	2.0
1	E	199	GLN	2.0

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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
2	MET	A	1001	9/9	0.99	0.04	17,21,23,23	0
2	MET	B	1001	9/9	0.99	0.05	16,19,24,24	0
2	MET	C	1001	9/9	0.99	0.04	17,20,22,23	0
2	MET	D	1001	9/9	0.99	0.05	17,21,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MET	E	1001	9/9	0.99	0.04	17,20,23,23	0
2	MET	F	1001	9/9	0.99	0.04	16,19,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.