



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 02:57 AM UTC

PDB ID : 3GQQ / pdb_00003gqq
Title : Crystal structure of the human retinal protein 4 (unc-119 homolog A). Northeast Structural Genomics Consortium target HR3066a
Authors : Vorobiev, S.M.; Chen, Y.; Seetharaman, J.; Shastry, R.; Foote, E.L.; Ciccosanti, C.; Sahdev, S.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2009-03-24
Resolution : 1.95 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:

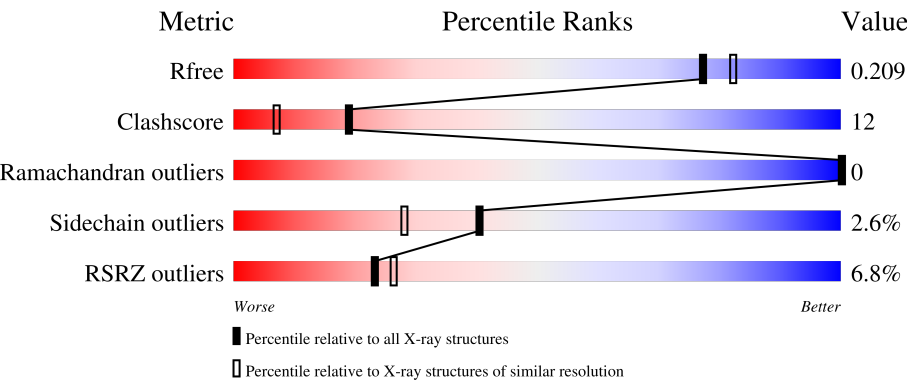
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 1.95 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}	180053	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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.

Mol	Chain	Length	Quality of chain
1	A	195	3% 68% 16% • 14%
1	B	195	7% 76% 14% • 9%
1	C	195	4% 67% 18% • 13%
1	D	195	7% 73% 19% • 7%

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Mol	Chain	Length	Quality of chain
1	E	195	6% 67% 15% • 15%
1	F	195	8% 67% 17% • 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9290 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 unc-119 homolog A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	Se	0	0	0
			1378	887	232	252	3	4			
1	B	178	Total	C	N	O	S	Se	0	0	0
			1458	936	247	268	3	4			
1	C	170	Total	C	N	O	S	Se	0	0	0
			1408	907	238	256	3	4			
1	D	181	Total	C	N	O	S	Se	0	0	0
			1483	953	252	271	3	4			
1	E	165	Total	C	N	O	S	Se	0	0	0
			1367	880	229	251	3	4			
1	F	164	Total	C	N	O	S	Se	0	0	0
			1363	879	229	248	3	4			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	MSE	-	expression tag	UNP Q13432
A	47	GLY	-	expression tag	UNP Q13432
A	48	HIS	-	expression tag	UNP Q13432
A	49	HIS	-	expression tag	UNP Q13432
A	50	HIS	-	expression tag	UNP Q13432
A	51	HIS	-	expression tag	UNP Q13432
A	52	HIS	-	expression tag	UNP Q13432
A	53	HIS	-	expression tag	UNP Q13432
A	54	SER	-	expression tag	UNP Q13432
A	55	HIS	-	expression tag	UNP Q13432
B	46	MSE	-	expression tag	UNP Q13432
B	47	GLY	-	expression tag	UNP Q13432
B	48	HIS	-	expression tag	UNP Q13432
B	49	HIS	-	expression tag	UNP Q13432
B	50	HIS	-	expression tag	UNP Q13432
B	51	HIS	-	expression tag	UNP Q13432
B	52	HIS	-	expression tag	UNP Q13432

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Chain	Residue	Modelled	Actual	Comment	Reference
B	53	HIS	-	expression tag	UNP Q13432
B	54	SER	-	expression tag	UNP Q13432
B	55	HIS	-	expression tag	UNP Q13432
C	46	MSE	-	expression tag	UNP Q13432
C	47	GLY	-	expression tag	UNP Q13432
C	48	HIS	-	expression tag	UNP Q13432
C	49	HIS	-	expression tag	UNP Q13432
C	50	HIS	-	expression tag	UNP Q13432
C	51	HIS	-	expression tag	UNP Q13432
C	52	HIS	-	expression tag	UNP Q13432
C	53	HIS	-	expression tag	UNP Q13432
C	54	SER	-	expression tag	UNP Q13432
C	55	HIS	-	expression tag	UNP Q13432
D	46	MSE	-	expression tag	UNP Q13432
D	47	GLY	-	expression tag	UNP Q13432
D	48	HIS	-	expression tag	UNP Q13432
D	49	HIS	-	expression tag	UNP Q13432
D	50	HIS	-	expression tag	UNP Q13432
D	51	HIS	-	expression tag	UNP Q13432
D	52	HIS	-	expression tag	UNP Q13432
D	53	HIS	-	expression tag	UNP Q13432
D	54	SER	-	expression tag	UNP Q13432
D	55	HIS	-	expression tag	UNP Q13432
E	46	MSE	-	expression tag	UNP Q13432
E	47	GLY	-	expression tag	UNP Q13432
E	48	HIS	-	expression tag	UNP Q13432
E	49	HIS	-	expression tag	UNP Q13432
E	50	HIS	-	expression tag	UNP Q13432
E	51	HIS	-	expression tag	UNP Q13432
E	52	HIS	-	expression tag	UNP Q13432
E	53	HIS	-	expression tag	UNP Q13432
E	54	SER	-	expression tag	UNP Q13432
E	55	HIS	-	expression tag	UNP Q13432
F	46	MSE	-	expression tag	UNP Q13432
F	47	GLY	-	expression tag	UNP Q13432
F	48	HIS	-	expression tag	UNP Q13432
F	49	HIS	-	expression tag	UNP Q13432
F	50	HIS	-	expression tag	UNP Q13432
F	51	HIS	-	expression tag	UNP Q13432
F	52	HIS	-	expression tag	UNP Q13432
F	53	HIS	-	expression tag	UNP Q13432
F	54	SER	-	expression tag	UNP Q13432

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Chain	Residue	Modelled	Actual	Comment	Reference
F	55	HIS	-	expression tag	UNP Q13432

- Molecule 2 is UNKNOWN LIGAND (CCD ID: UNL) (formula:).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	6	Total O 6 6	0	0
2	B	4	Total O 4 4	0	0
2	C	4	Total O 4 4	0	0
2	D	6	Total O 6 6	0	0
2	E	6	Total O 6 6	0	0
2	F	4	Total O 4 4	0	0

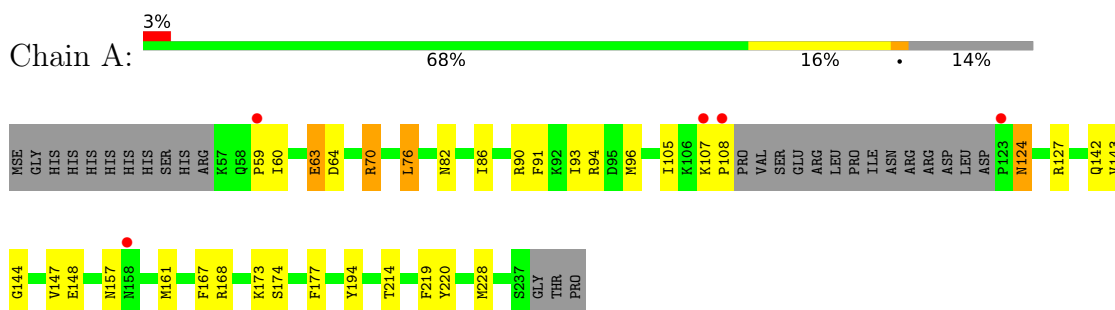
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	160	Total O 160 160	0	0
3	B	134	Total O 134 134	0	0
3	C	155	Total O 155 155	0	0
3	D	158	Total O 158 158	0	0
3	E	108	Total O 108 108	0	0
3	F	88	Total O 88 88	0	0

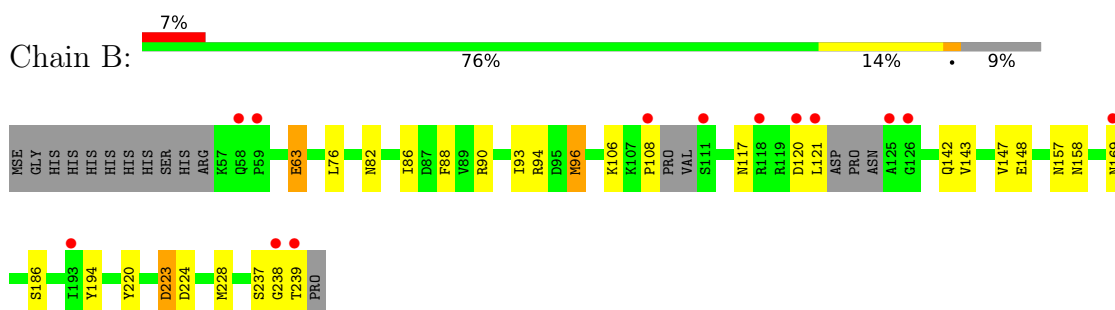
3 Residue-property plots [i](#)

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.

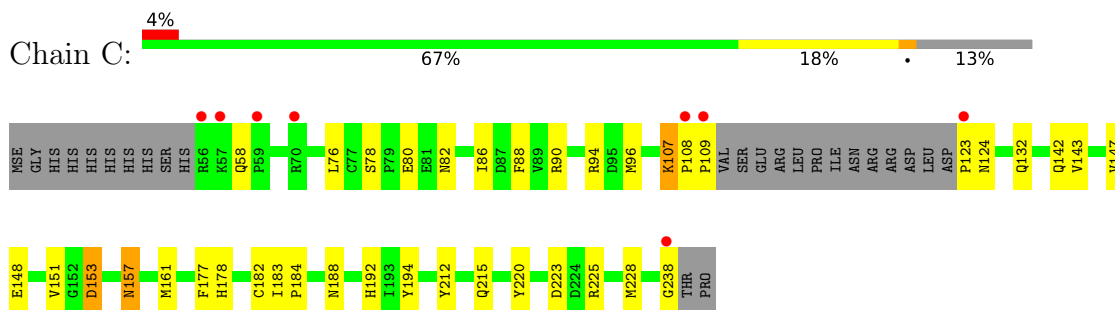
• Molecule 1: Protein unc-119 homolog A



• Molecule 1: Protein unc-119 homolog A

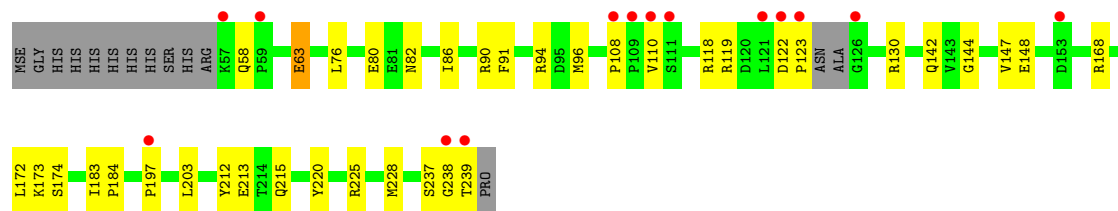


• Molecule 1: Protein unc-119 homolog A

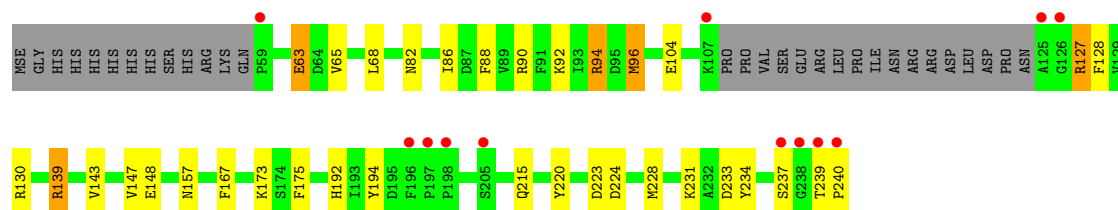


• Molecule 1: Protein unc-119 homolog A

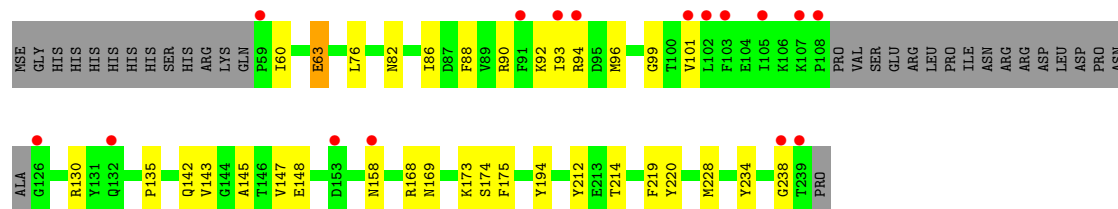




● Molecule 1: Protein unc-119 homolog A



● Molecule 1: Protein unc-119 homolog A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.89Å 79.56Å 189.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.51 – 1.95 40.51 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.0 (40.51-1.95) 99.3 (40.51-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 1.94Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.191 , 0.214 0.188 , 0.209	Depositor DCC
R_{free} test set	4351 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	25.8	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9290	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.35 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.8646e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: UNL

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	0/1412	0.87	3/1900 (0.2%)
1	B	0.34	0/1492	0.85	1/2007 (0.0%)
1	C	0.35	0/1444	0.82	1/1942 (0.1%)
1	D	0.36	0/1520	0.87	2/2050 (0.1%)
1	E	0.33	0/1401	0.84	2/1883 (0.1%)
1	F	0.32	0/1397	0.81	2/1877 (0.1%)
All	All	0.35	0/8666	0.84	11/11659 (0.1%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	124	ASN	N-CA-C	-7.49	103.13	111.82
1	E	127	ARG	N-CA-C	-7.31	104.51	113.50
1	A	124	ASN	N-CA-C	-6.77	104.28	112.54
1	A	108	PRO	N-CA-CB	6.38	110.02	103.00
1	F	60	ILE	N-CA-C	5.51	116.52	108.58
1	A	214	THR	N-CA-C	-5.28	100.13	108.73
1	B	238	GLY	N-CA-C	-5.26	106.78	114.61
1	F	214	THR	N-CA-C	-5.15	100.34	108.73
1	E	167	PHE	N-CA-C	-5.04	100.20	108.41
1	D	58	GLN	CA-C-N	5.02	125.00	120.03
1	D	58	GLN	C-N-CA	5.02	125.00	120.03

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	A	1378	0	1310	30	0
1	B	1458	0	1386	27	0
1	C	1408	0	1356	48	0
1	D	1483	0	1414	34	0
1	E	1367	0	1309	36	0
1	F	1363	0	1310	26	0
2	A	6	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	6	0	0	1	0
2	E	6	0	0	0	0
2	F	4	0	0	0	0
3	A	160	0	0	3	0
3	B	134	0	0	5	0
3	C	155	0	0	4	0
3	D	158	0	0	1	0
3	E	108	0	0	4	0
3	F	88	0	0	2	0
All	All	9290	0	8085	196	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 (196) 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:96:MSE:HA	1:E:96:MSE:HE2	1.28	1.08
1:B:96:MSE:HA	1:B:96:MSE:HE2	1.35	1.07
1:E:239:THR:N	1:E:240:PRO:HA	1.75	1.01
1:E:139:ARG:HH11	1:E:139:ARG:HG3	1.26	1.00
1:C:183:ILE:H	1:C:188:ASN:HD21	1.03	0.99
1:C:107:LYS:HB3	1:C:109:PRO:HB3	1.45	0.97
1:C:96:MSE:HA	1:C:96:MSE:HE2	1.47	0.94
1:C:238:GLY:HA2	1:D:238:GLY:HA3	1.49	0.92
1:E:139:ARG:HH11	1:E:139:ARG:CG	1.85	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:MSE:HE3	1:D:144:GLY:N	1.86	0.89
1:E:96:MSE:HA	1:E:96:MSE:CE	2.01	0.89
1:D:96:MSE:HE3	1:D:144:GLY:CA	2.02	0.89
1:A:96:MSE:HE3	1:A:144:GLY:H	1.37	0.88
1:A:96:MSE:HE3	1:A:144:GLY:N	1.89	0.88
1:D:96:MSE:HE3	1:D:144:GLY:HA3	1.59	0.84
1:B:96:MSE:HA	1:B:96:MSE:CE	2.08	0.84
1:E:63:GLU:H	1:E:63:GLU:CD	1.83	0.84
1:E:94:ARG:HH11	1:E:96:MSE:HE1	1.40	0.84
1:C:215:GLN:HE22	1:E:157:ASN:HD21	1.26	0.84
1:E:239:THR:H	1:E:240:PRO:HA	1.39	0.83
1:D:63:GLU:H	1:D:63:GLU:CD	1.87	0.81
1:B:106:LYS:O	1:B:108:PRO:HD3	1.81	0.81
1:A:70:ARG:HG2	1:A:70:ARG:HH11	1.46	0.80
1:D:96:MSE:HE3	1:D:144:GLY:H	1.49	0.76
1:C:94:ARG:HG3	1:C:96:MSE:CE	2.16	0.75
1:D:94:ARG:O	1:D:96:MSE:HE2	1.87	0.75
1:C:108:PRO:N	1:C:109:PRO:HA	2.03	0.74
1:C:94:ARG:HG3	1:C:96:MSE:HE1	1.70	0.73
1:E:94:ARG:NH1	1:E:96:MSE:HE1	2.03	0.73
1:F:96:MSE:HG3	1:F:142:GLN:HG2	1.69	0.73
1:E:239:THR:N	1:E:240:PRO:CA	2.51	0.72
1:F:94:ARG:HG3	1:F:96:MSE:CE	2.19	0.72
1:B:63:GLU:CD	1:B:63:GLU:H	1.97	0.71
1:C:215:GLN:HE22	1:E:157:ASN:ND2	1.91	0.69
1:F:90:ARG:HB3	1:F:148:GLU:HB2	1.74	0.69
1:A:96:MSE:HE3	1:A:144:GLY:CA	2.23	0.68
1:C:94:ARG:NE	1:C:96:MSE:HE1	2.09	0.68
1:F:94:ARG:HH21	1:F:99:GLY:HA2	1.59	0.68
1:A:94:ARG:O	1:A:96:MSE:HE2	1.94	0.66
1:B:94:ARG:HG3	1:B:96:MSE:HE3	1.76	0.66
1:A:157:ASN:HD21	1:D:215:GLN:HE22	1.46	0.64
1:C:96:MSE:HE2	1:C:96:MSE:CA	2.25	0.64
1:D:168:ARG:NH1	1:D:213:GLU:OE1	2.31	0.63
1:B:94:ARG:HG3	1:B:96:MSE:CE	2.28	0.63
1:C:94:ARG:HE	1:C:96:MSE:HE1	1.62	0.63
1:B:90:ARG:HB3	1:B:148:GLU:HB2	1.81	0.62
1:C:183:ILE:N	1:C:188:ASN:HD21	1.86	0.62
1:C:96:MSE:HA	1:C:96:MSE:CE	2.27	0.62
1:C:108:PRO:CD	1:C:109:PRO:HA	2.30	0.62
1:A:94:ARG:HB3	1:A:96:MSE:HE1	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:130:ARG:HG3	1:E:233:ASP:OD1	2.01	0.60
1:A:70:ARG:HG2	1:A:70:ARG:NH1	2.13	0.59
1:A:63:GLU:H	1:A:63:GLU:CD	2.11	0.59
1:E:139:ARG:CG	1:E:139:ARG:NH1	2.54	0.59
1:B:117:ASN:O	1:B:120:ASP:HB2	2.03	0.59
1:F:169:ASN:HA	3:F:717:HOH:O	2.02	0.58
1:C:238:GLY:CA	1:D:238:GLY:HA3	2.30	0.57
1:A:82:ASN:HD21	1:A:86:ILE:H	1.51	0.57
1:C:90:ARG:HB3	1:C:148:GLU:HB2	1.86	0.57
1:C:107:LYS:HB3	1:C:109:PRO:CB	2.29	0.57
1:C:183:ILE:H	1:C:188:ASN:ND2	1.87	0.57
1:F:130:ARG:HD3	1:F:212:TYR:CE1	2.39	0.57
1:B:94:ARG:NE	1:B:96:MSE:HE1	2.20	0.56
1:B:88:PHE:CD2	1:B:147:VAL:HG21	2.41	0.56
1:C:94:ARG:O	1:C:96:MSE:HE3	2.06	0.55
1:C:94:ARG:HG3	1:C:96:MSE:HE3	1.88	0.55
1:E:82:ASN:HD21	1:E:86:ILE:H	1.53	0.55
1:E:94:ARG:CG	1:E:96:MSE:HE3	2.37	0.55
1:F:94:ARG:HD3	3:F:1175:HOH:O	2.06	0.55
1:C:96:MSE:HG3	1:C:142:GLN:HE21	1.72	0.55
1:A:96:MSE:HE3	1:A:144:GLY:HA3	1.89	0.54
1:E:94:ARG:HG2	1:E:96:MSE:HE3	1.88	0.54
1:C:178:HIS:HD2	3:C:674:HOH:O	1.90	0.54
1:C:108:PRO:HG2	1:C:109:PRO:O	2.08	0.54
1:C:82:ASN:HD21	1:C:86:ILE:H	1.54	0.54
1:D:96:MSE:HG3	1:D:142:GLN:HE21	1.73	0.54
1:B:82:ASN:HD21	1:B:86:ILE:H	1.57	0.53
1:C:108:PRO:HG2	1:C:109:PRO:C	2.33	0.53
1:D:118:ARG:HG2	1:D:119:ARG:HG3	1.90	0.53
1:E:127:ARG:NH1	3:E:657:HOH:O	2.41	0.53
1:C:88:PHE:HD2	1:C:147:VAL:CG2	2.21	0.53
1:F:82:ASN:HD21	1:F:86:ILE:H	1.56	0.52
1:A:90:ARG:HB3	1:A:148:GLU:HB2	1.91	0.52
1:B:157:ASN:HD22	1:B:223:ASP:CG	2.18	0.52
1:D:96:MSE:CE	1:D:144:GLY:H	2.18	0.52
1:F:173:LYS:HG2	1:F:175:PHE:CZ	2.45	0.52
1:C:215:GLN:NE2	1:E:157:ASN:HD21	2.02	0.52
1:A:173:LYS:HG3	1:A:174:SER:N	2.25	0.51
1:C:220:TYR:HB2	1:C:228:MSE:HB2	1.91	0.51
1:A:94:ARG:HB3	1:A:96:MSE:CE	2.41	0.51
1:D:63:GLU:CD	1:D:63:GLU:N	2.65	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:173:LYS:HD3	1:D:197:PRO:HD3	1.92	0.51
1:D:220:TYR:HB2	1:D:228:MSE:HB2	1.92	0.51
1:C:132:GLN:NE2	1:C:212:TYR:OH	2.35	0.50
1:D:94:ARG:HB3	1:D:96:MSE:HE1	1.93	0.50
1:B:88:PHE:CD2	1:B:147:VAL:CG2	2.94	0.50
1:C:153:ASP:OD2	1:C:153:ASP:N	2.44	0.50
1:F:168:ARG:O	1:F:169:ASN:HB2	2.12	0.50
1:D:82:ASN:HD21	1:D:86:ILE:H	1.58	0.50
1:D:238:GLY:O	1:D:239:THR:CB	2.58	0.50
1:F:173:LYS:HG3	1:F:174:SER:N	2.26	0.50
1:C:94:ARG:CG	1:C:96:MSE:HE1	2.40	0.50
1:C:143:VAL:HG22	1:C:194:TYR:HB2	1.93	0.49
1:D:173:LYS:HG3	1:D:174:SER:N	2.27	0.49
1:A:91:PHE:HD1	1:A:147:VAL:HG12	1.77	0.49
1:C:151:VAL:HG23	1:C:184:PRO:HA	1.95	0.49
1:C:123:PRO:HG2	3:C:961:HOH:O	2.11	0.49
1:E:215:GLN:HG2	1:E:233:ASP:HB3	1.94	0.49
1:D:90:ARG:HB3	1:D:148:GLU:HB2	1.94	0.49
1:A:167:PHE:CE2	1:A:168:ARG:HD2	2.48	0.48
1:E:143:VAL:CG2	1:E:194:TYR:HB2	2.43	0.48
1:C:88:PHE:CD2	1:C:147:VAL:CG2	2.96	0.48
1:E:139:ARG:NH1	3:E:1282:HOH:O	2.45	0.48
1:C:108:PRO:N	1:C:109:PRO:CA	2.75	0.48
1:E:63:GLU:CD	1:E:63:GLU:N	2.62	0.48
1:F:220:TYR:HB2	1:F:228:MSE:HB2	1.95	0.47
1:B:220:TYR:HB2	1:B:228:MSE:HB2	1.96	0.47
1:E:90:ARG:HB3	1:E:148:GLU:HB2	1.95	0.47
1:F:94:ARG:HG3	1:F:96:MSE:HE3	1.95	0.47
1:E:96:MSE:CE	1:E:96:MSE:CA	2.85	0.47
1:E:139:ARG:HG3	1:E:139:ARG:NH1	2.08	0.47
1:A:94:ARG:O	1:A:96:MSE:CE	2.62	0.46
1:B:88:PHE:CE2	1:B:147:VAL:HG21	2.51	0.46
1:F:82:ASN:C	1:F:82:ASN:HD22	2.23	0.46
1:F:143:VAL:HG22	1:F:194:TYR:HB2	1.97	0.46
1:C:225:ARG:NE	3:C:715:HOH:O	2.48	0.46
1:D:168:ARG:NH1	3:D:960:HOH:O	2.48	0.46
1:C:108:PRO:CG	1:C:109:PRO:HA	2.45	0.46
1:D:172:LEU:HD22	1:D:203:LEU:HD21	1.98	0.46
1:D:94:ARG:HB3	1:D:96:MSE:CE	2.46	0.46
1:A:220:TYR:HB2	1:A:228:MSE:HB2	1.98	0.46
1:C:182:CYS:O	1:C:184:PRO:HD3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:223:ASP:O	1:E:224:ASP:HB2	2.15	0.46
1:A:124:ASN:O	1:A:127:ARG:HG3	2.16	0.45
1:F:92:LYS:HE2	1:F:101:VAL:HG22	1.98	0.45
1:F:96:MSE:HG3	1:F:142:GLN:HE21	1.81	0.45
1:A:105:ILE:HD11	3:A:967:HOH:O	2.17	0.45
1:D:225:ARG:HH11	1:D:225:ARG:HG3	1.82	0.45
1:A:124:ASN:O	1:A:127:ARG:CG	2.64	0.45
1:B:96:MSE:HG3	1:B:142:GLN:HE21	1.82	0.45
1:A:161:MSE:HE2	1:A:177:PHE:HD2	1.81	0.44
1:D:122:ASP:HA	1:D:123:PRO:HD3	1.84	0.44
1:F:234:TYR:CD1	1:F:234:TYR:N	2.85	0.44
1:B:88:PHE:HD2	1:B:147:VAL:CG2	2.30	0.44
1:B:96:MSE:CE	1:B:96:MSE:CA	2.88	0.44
1:E:128:PHE:HA	1:E:231:LYS:O	2.17	0.44
1:A:96:MSE:HG3	1:A:142:GLN:HE21	1.83	0.44
1:F:88:PHE:HD2	1:F:147:VAL:HG22	1.81	0.44
1:D:118:ARG:O	1:D:119:ARG:HB2	2.18	0.44
1:D:91:PHE:HD1	1:D:147:VAL:HG12	1.83	0.44
1:C:88:PHE:CD2	1:C:147:VAL:HG21	2.53	0.43
1:E:234:TYR:CD1	1:E:234:TYR:N	2.85	0.43
1:B:143:VAL:CG2	1:B:194:TYR:HB2	2.48	0.43
1:A:60:ILE:HG23	1:A:64:ASP:HB2	1.99	0.43
1:A:94:ARG:C	1:A:96:MSE:HE2	2.42	0.43
1:F:93:ILE:HG22	1:F:145:ALA:HB2	2.00	0.43
1:B:237:SER:C	1:B:239:THR:N	2.76	0.43
1:B:186:SER:HB2	3:B:742:HOH:O	2.19	0.43
1:D:80:GLU:OE1	1:D:80:GLU:N	2.45	0.43
1:F:63:GLU:CD	1:F:63:GLU:H	2.26	0.43
1:B:157:ASN:HB2	3:B:868:HOH:O	2.19	0.43
1:F:90:ARG:HD3	1:F:148:GLU:OE2	2.19	0.43
1:A:59:PRO:HB2	3:A:1190:HOH:O	2.19	0.43
1:E:228:MSE:HE3	3:E:1114:HOH:O	2.19	0.43
1:C:108:PRO:HG2	1:C:109:PRO:HA	2.01	0.43
1:B:120:ASP:O	1:B:121:LEU:C	2.62	0.42
1:B:169:ASN:HA	3:B:581:HOH:O	2.19	0.42
1:F:94:ARG:NH2	1:F:99:GLY:HA2	2.31	0.42
1:E:92:LYS:HG3	1:E:104:GLU:HB2	2.01	0.42
1:C:157:ASN:HB3	1:C:223:ASP:OD1	2.18	0.42
1:A:91:PHE:HB2	3:A:1134:HOH:O	2.18	0.42
1:F:76:LEU:HD22	1:F:219:PHE:CE2	2.55	0.42
1:E:88:PHE:HD2	1:E:147:VAL:HG22	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:MSE:HE2	1:C:177:PHE:HD2	1.85	0.42
1:B:158:ASN:ND2	3:B:1009:HOH:O	2.38	0.42
1:E:173:LYS:HG2	1:E:175:PHE:CZ	2.55	0.42
1:E:192:HIS:HE1	3:E:912:HOH:O	2.03	0.42
1:D:130:ARG:HD3	1:D:212:TYR:CE1	2.56	0.41
1:B:223:ASP:O	1:B:224:ASP:HB2	2.19	0.41
1:C:108:PRO:HG2	1:C:109:PRO:CA	2.51	0.41
1:D:183:ILE:HA	1:D:184:PRO:HD3	1.96	0.41
1:A:76:LEU:HD22	1:A:219:PHE:CE2	2.56	0.41
1:B:239:THR:CB	3:B:565:HOH:O	2.68	0.41
1:C:94:ARG:CD	1:C:96:MSE:HE1	2.51	0.41
1:D:96:MSE:CE	1:D:144:GLY:HA3	2.42	0.41
1:D:108:PRO:HG2	1:D:110:VAL:O	2.20	0.41
1:E:220:TYR:HB2	1:E:228:MSE:HB2	2.03	0.41
1:C:78:SER:HB2	1:C:80:GLU:OE1	2.21	0.40
1:E:65:VAL:HA	1:E:68:LEU:HG	2.02	0.40
1:A:96:MSE:CE	1:A:144:GLY:H	2.20	0.40
3:C:1081:HOH:O	1:D:237:SER:HB3	2.21	0.40
1:F:92:LYS:HE2	1:F:101:VAL:CG2	2.50	0.40
2:D:431:UNL:OH2	2:D:432:UNL:OH2	2.40	0.40
1:A:143:VAL:HG22	1:A:194:TYR:HB2	2.02	0.40
1:C:178:HIS:O	1:C:192:HIS:HE1	2.05	0.40
1:F:135:PRO:HD3	1:F:238:GLY:O	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	163/195 (84%)	163 (100%)	0	0	100	100
1	B	172/195 (88%)	166 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	166/195 (85%)	165 (99%)	1 (1%)	0	100	100
1	D	177/195 (91%)	174 (98%)	3 (2%)	0	100	100
1	E	161/195 (83%)	158 (98%)	3 (2%)	0	100	100
1	F	160/195 (82%)	159 (99%)	1 (1%)	0	100	100
All	All	999/1170 (85%)	985 (99%)	14 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	151/176 (86%)	146 (97%)	5 (3%)	33	20
1	B	159/176 (90%)	154 (97%)	5 (3%)	35	21
1	C	156/176 (89%)	151 (97%)	5 (3%)	34	20
1	D	163/176 (93%)	161 (99%)	2 (1%)	63	57
1	E	151/176 (86%)	146 (97%)	5 (3%)	33	20
1	F	151/176 (86%)	149 (99%)	2 (1%)	61	54
All	All	931/1056 (88%)	907 (97%)	24 (3%)	40	28

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	GLU
1	A	70	ARG
1	A	76	LEU
1	A	93	ILE
1	A	107	LYS
1	B	63	GLU
1	B	76	LEU
1	B	93	ILE
1	B	96	MSE

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Mol	Chain	Res	Type
1	B	223	ASP
1	C	58	GLN
1	C	76	LEU
1	C	107	LYS
1	C	153	ASP
1	C	157	ASN
1	D	63	GLU
1	D	76	LEU
1	E	63	GLU
1	E	94	ARG
1	E	96	MSE
1	E	139	ARG
1	E	237	SER
1	F	63	GLU
1	F	158	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	ASN
1	A	142	GLN
1	A	157	ASN
1	B	82	ASN
1	B	117	ASN
1	B	142	GLN
1	B	157	ASN
1	B	178	HIS
1	B	192	HIS
1	C	58	GLN
1	C	82	ASN
1	C	142	GLN
1	C	178	HIS
1	C	188	ASN
1	C	192	HIS
1	D	82	ASN
1	D	142	GLN
1	D	157	ASN
1	D	192	HIS
1	E	82	ASN
1	E	157	ASN
1	E	158	ASN
1	E	192	HIS

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Mol	Chain	Res	Type
1	E	210	HIS
1	F	82	ASN
1	F	142	GLN
1	F	157	ASN
1	F	210	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 30 are unknown - leaving 0 for Mogul analysis.

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.

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	163/195 (83%)	-0.04	5 (3%)	51	57	14, 24, 42, 69	0
1	B	174/195 (89%)	0.30	13 (7%)	20	23	16, 31, 74, 90	0
1	C	166/195 (85%)	0.28	8 (4%)	35	40	15, 29, 55, 102	0
1	D	177/195 (90%)	0.09	14 (7%)	18	21	13, 26, 56, 83	0
1	E	161/195 (82%)	0.36	12 (7%)	20	23	16, 33, 61, 107	0
1	F	160/195 (82%)	0.62	16 (10%)	12	14	18, 41, 68, 102	0
All	All	1001/1170 (85%)	0.27	68 (6%)	23	26	13, 30, 64, 107	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	238	GLY	7.4
1	F	239	THR	6.2
1	C	56	ARG	5.6
1	C	109	PRO	5.3
1	E	239	THR	4.8
1	E	240	PRO	4.7
1	E	238	GLY	4.6
1	D	110	VAL	4.6
1	E	125	ALA	4.5
1	B	239	THR	4.4
1	D	239	THR	4.2
1	C	108	PRO	4.2
1	A	59	PRO	4.1
1	B	108	PRO	3.8
1	B	125	ALA	3.8
1	D	123	PRO	3.7
1	B	121	LEU	3.5
1	F	103	PHE	3.3
1	D	109	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	E	197	PRO	3.2
1	A	123	PRO	3.1
1	F	238	GLY	3.1
1	B	126	GLY	2.9
1	F	126	GLY	2.9
1	C	59	PRO	2.9
1	B	118	ARG	2.9
1	B	238	GLY	2.8
1	D	108	PRO	2.8
1	E	196	PHE	2.8
1	F	153	ASP	2.8
1	E	126	GLY	2.7
1	F	102	LEU	2.7
1	E	59	PRO	2.7
1	E	107	LYS	2.6
1	B	59	PRO	2.6
1	B	120	ASP	2.6
1	F	59	PRO	2.5
1	D	111	SER	2.5
1	C	70	ARG	2.5
1	D	122	ASP	2.5
1	F	93	ILE	2.4
1	E	198	PRO	2.4
1	C	123	PRO	2.3
1	B	111	SER	2.3
1	F	158	ASN	2.3
1	E	205	SER	2.3
1	F	107	LYS	2.3
1	A	108	PRO	2.3
1	D	238	GLY	2.3
1	D	197	PRO	2.2
1	A	158	ASN	2.2
1	D	57	LYS	2.2
1	F	91	PHE	2.2
1	D	126	GLY	2.2
1	F	108	PRO	2.2
1	F	105	ILE	2.1
1	E	237	SER	2.1
1	A	107	LYS	2.1
1	C	57	LYS	2.1
1	B	58	GLN	2.1
1	F	101	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	153	ASP	2.1
1	F	94	ARG	2.1
1	B	193	ILE	2.0
1	F	132	GLN	2.0
1	D	121	LEU	2.0
1	B	169	ASN	2.0
1	D	59	PRO	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(Å ²)	Q<0.9
2	UNL	F	452	1/-	0.71	0.25	66,66,66,66	0
2	UNL	B	411	1/-	0.77	0.32	66,66,66,66	0
2	UNL	A	401	1/-	0.77	0.38	60,60,60,60	0
2	UNL	F	453	1/-	0.80	0.18	58,58,58,58	0
2	UNL	C	421	1/-	0.82	0.37	55,55,55,55	0
2	UNL	A	405	1/-	0.84	0.28	53,53,53,53	0
2	UNL	D	436	1/-	0.85	0.21	52,52,52,52	0
2	UNL	F	451	1/-	0.85	0.30	62,62,62,62	0
2	UNL	D	431	1/-	0.85	0.31	54,54,54,54	0
2	UNL	D	433	1/-	0.85	0.16	52,52,52,52	0
2	UNL	B	410	1/-	0.86	0.31	59,59,59,59	0
2	UNL	C	423	1/-	0.86	0.24	49,49,49,49	0
2	UNL	E	442	1/-	0.87	0.31	57,57,57,57	0
2	UNL	D	435	1/-	0.90	0.22	38,38,38,38	0
2	UNL	E	441	1/-	0.90	0.28	53,53,53,53	0
2	UNL	A	404	1/-	0.91	0.16	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	UNL	C	422	1/-	0.91	0.33	54,54,54,54	0
2	UNL	A	403	1/-	0.91	0.25	48,48,48,48	0
2	UNL	B	412	1/-	0.91	0.23	46,46,46,46	0
2	UNL	D	432	1/-	0.92	0.20	52,52,52,52	0
2	UNL	B	413	1/-	0.92	0.21	47,47,47,47	0
2	UNL	E	444	1/-	0.92	0.09	50,50,50,50	0
2	UNL	A	402	1/-	0.93	0.23	36,36,36,36	0
2	UNL	E	443	1/-	0.93	0.11	46,46,46,46	0
2	UNL	F	454	1/-	0.93	0.09	37,37,37,37	0
2	UNL	C	424	1/-	0.94	0.07	31,31,31,31	0
2	UNL	D	434	1/-	0.95	0.19	40,40,40,40	0
2	UNL	E	445	1/-	0.95	0.14	31,31,31,31	0
2	UNL	A	406	1/-	0.95	0.13	32,32,32,32	0
2	UNL	E	446	1/-	0.99	0.11	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.