



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 08:47 pm BST

PDB ID : 6U9S
Title : Crystal structure of human CD81 large extracellular loop in complex with 5A6 Fab
Authors : Susa, K.J.; Seegar, T.C.M.; Blacklow, S.C.B.; Kruse, A.C.
Deposited on : 2019-09-09
Resolution : 2.40 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 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) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

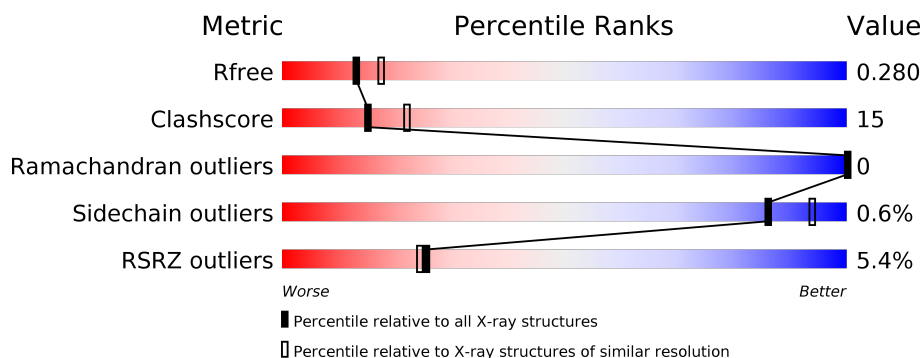
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 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	232	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div>5%</div> </div> </div>
1	D	232	<div> <div>15%</div> <div> <div></div> <div>68%</div> <div>26%</div> <div>6%</div> </div> </div>
2	B	219	<div> <div>70%</div> <div>30%</div> </div>
2	E	219	<div> <div>9%</div> <div> <div></div> <div>69%</div> <div>30%</div> </div> </div>
3	C	98	<div> <div>69%</div> <div>23%</div> <div>• 6%</div> </div>
3	F	98	<div> <div>%</div> <div> <div></div> <div>69%</div> <div>23%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8374 atoms, of which 24 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 5A6 FAB Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1641	1044	265	327	5			
1	D	218	Total	C	N	O	S	0	0	0
			1593	1012	255	321	5			

- Molecule 2 is a protein called 5A6 FAB Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	219	Total	C	N	O	S	0	0	0
			1674	1051	279	338	6			
2	E	218	Total	C	N	O	S	0	0	0
			1666	1045	279	337	5			

- Molecule 3 is a protein called CD81 antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	92	Total	C	N	O	S	0	0	0
			701	433	119	145	4			
3	F	92	Total	C	N	O	S	0	0	0
			699	432	119	144	4			

There are 18 discrepancies between the modelled and reference sequences:

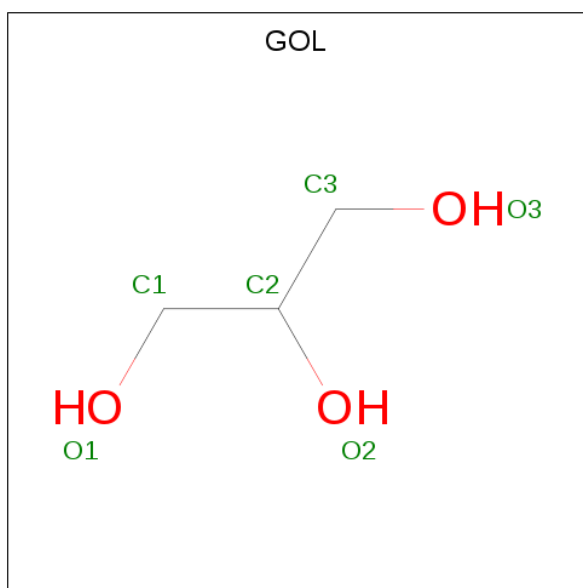
Chain	Residue	Modelled	Actual	Comment	Reference
C	110	GLU	-	expression tag	UNP P60033
C	111	THR	-	expression tag	UNP P60033
C	201	SER	-	expression tag	UNP P60033
C	202	HIS	-	expression tag	UNP P60033
C	203	HIS	-	expression tag	UNP P60033
C	204	HIS	-	expression tag	UNP P60033
C	205	HIS	-	expression tag	UNP P60033
C	206	HIS	-	expression tag	UNP P60033

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Chain	Residue	Modelled	Actual	Comment	Reference
C	207	HIS	-	expression tag	UNP P60033
F	110	GLU	-	expression tag	UNP P60033
F	111	THR	-	expression tag	UNP P60033
F	201	SER	-	expression tag	UNP P60033
F	202	HIS	-	expression tag	UNP P60033
F	203	HIS	-	expression tag	UNP P60033
F	204	HIS	-	expression tag	UNP P60033
F	205	HIS	-	expression tag	UNP P60033
F	206	HIS	-	expression tag	UNP P60033
F	207	HIS	-	expression tag	UNP P60033

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	E	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	77	Total	O	0	0
			77	77		

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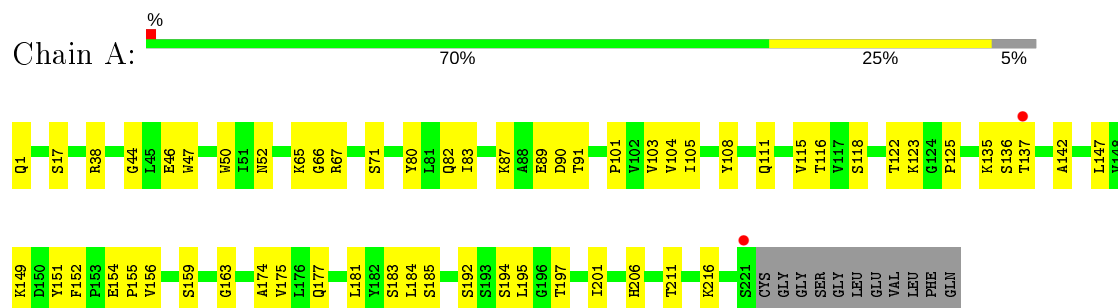
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	94	Total 94	O 94	0	0
5	C	69	Total 69	O 69	0	0
5	D	25	Total 25	O 25	0	0
5	E	40	Total 40	O 40	0	0
5	F	53	Total 53	O 53	0	0

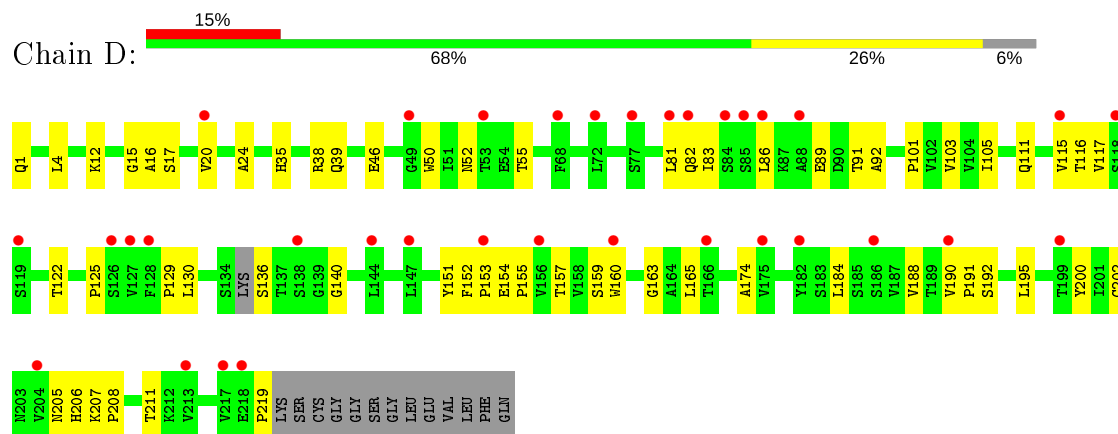
3 Residue-property plots [i](#)

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.

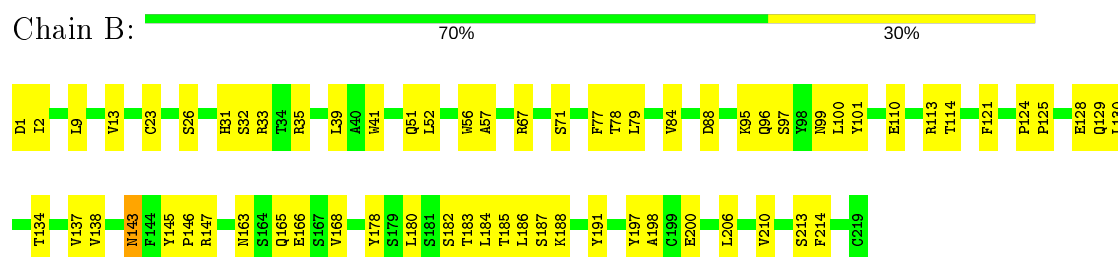
• Molecule 1: 5A6 FAB Heavy Chain



• Molecule 1: 5A6 FAB Heavy Chain

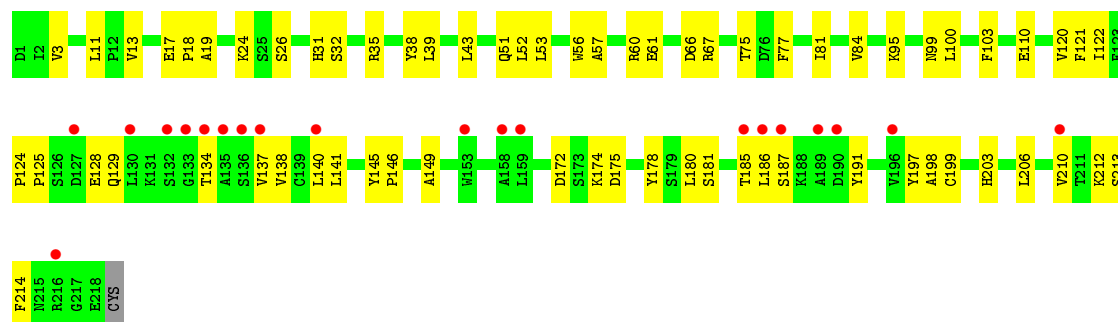


• Molecule 2: 5A6 FAB Light Chain



• Molecule 2: 5A6 FAB Light Chain





- Molecule 3: CD81 antigen

Chain C: 69% 23% 6%



- Molecule 3: CD81 antigen

Chain F: 69% 23% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	40.00Å 96.86Å 297.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.52 – 2.40 49.52 – 2.40	Depositor EDS
% Data completeness (in resolution range)	87.0 (49.52-2.40) 84.9 (49.52-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.67 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.16_3549, PHENIX 1.16_3549	Depositor
R, R_{free}	0.214 , 0.281 0.214 , 0.280	Depositor DCC
R_{free} test set	2000 reflections (4.31%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8374	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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](#)

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

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	0.27	0/1683	0.47	0/2304
1	D	0.26	0/1634	0.47	0/2246
2	B	0.26	0/1712	0.46	0/2330
2	E	0.25	0/1704	0.43	0/2321
3	C	0.26	0/710	0.43	0/960
3	F	0.26	0/708	0.43	0/957
All	All	0.26	0/8151	0.45	0/11118

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	1641	0	1591	53	0
1	D	1593	0	1500	53	0
2	B	1674	0	1604	51	0
2	E	1666	0	1588	55	0
3	C	701	0	674	19	0
3	F	699	0	669	21	0
4	A	6	8	8	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	6	8	8	0	0
4	E	6	8	8	0	0
5	A	77	0	0	10	1
5	B	94	0	0	8	0
5	C	69	0	0	5	3
5	D	25	0	0	1	0
5	E	40	0	0	2	0
5	F	53	0	0	4	2
All	All	8350	24	7650	235	3

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 15.

All (235) 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:D:91:THR:HG23	1:D:116:THR:HA	1.50	0.91
2:B:51:GLN:NE2	5:B:401:HOH:O	2.05	0.87
1:A:66:GLY:HA2	3:F:162:LEU:HD22	1.57	0.85
1:D:125:PRO:HB3	1:D:151:TYR:HB3	1.59	0.84
2:E:191:TYR:OH	5:E:401:HOH:O	1.96	0.83
1:A:192:SER:HA	1:A:195:LEU:HD13	1.62	0.81
1:D:86:LEU:HG	1:D:117:VAL:HG21	1.64	0.79
3:F:132:GLN:OE1	5:F:301:HOH:O	1.98	0.79
3:F:133:GLN:OE1	5:F:302:HOH:O	1.99	0.77
3:C:135:VAL:HG21	3:C:165:LEU:HD22	1.66	0.77
3:C:137:ASP:OD2	3:C:139:ASP:N	2.14	0.77
1:D:206:HIS:CE1	1:D:208:PRO:HG2	2.21	0.76
1:A:185:SER:HB3	5:A:559:HOH:O	1.86	0.73
1:A:91:THR:HG23	1:A:116:THR:HA	1.72	0.71
1:D:105:ILE:HD12	2:E:52:LEU:CD2	2.20	0.71
1:D:130:LEU:HD21	2:E:138:VAL:HG21	1.74	0.68
1:A:192:SER:HA	1:A:195:LEU:CD1	2.24	0.67
2:E:32:SER:HA	2:E:35:ARG:HH21	1.59	0.67
3:F:184:ASN:HA	3:F:187:LYS:HE2	1.76	0.67
1:A:105:ILE:HD12	2:B:52:LEU:CD2	2.25	0.66
2:E:206:LEU:O	5:E:402:HOH:O	2.14	0.65
2:E:198:ALA:HB2	2:E:213:SER:CB	2.26	0.65
3:F:133:GLN:NE2	5:F:304:HOH:O	2.19	0.65
2:B:129:GLN:NE2	5:B:402:HOH:O	2.11	0.65
3:F:132:GLN:O	3:F:136:VAL:HG13	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:111:THR:HG22	3:C:119:ILE:HD11	1.81	0.63
2:B:206:LEU:HD13	2:B:210:VAL:HG23	1.81	0.63
3:F:141:ASN:O	5:F:303:HOH:O	2.16	0.62
2:B:191:TYR:O	2:B:197:TYR:OH	2.18	0.62
2:B:57:ALA:HA	5:B:465:HOH:O	2.01	0.61
2:E:198:ALA:HB2	2:E:213:SER:HB3	1.81	0.61
1:D:52:ASN:ND2	3:F:180:ASN:HB3	2.15	0.61
1:D:82:GLN:HB3	5:D:317:HOH:O	2.01	0.61
2:B:110:GLU:OE1	2:B:178:TYR:OH	2.17	0.61
1:D:206:HIS:HB3	1:D:211:THR:OG1	2.00	0.61
1:A:71:SER:OG	1:A:80:TYR:HB2	2.01	0.60
2:E:39:LEU:HD22	2:E:77:PHE:CD1	2.36	0.60
2:E:13:VAL:HG11	2:E:84:VAL:HG21	1.83	0.60
1:D:105:ILE:HD12	2:E:52:LEU:HD22	1.83	0.60
2:E:128:GLU:OE1	2:E:128:GLU:N	2.35	0.60
1:D:140:GLY:O	1:D:192:SER:N	2.33	0.59
2:E:43:LEU:HB2	2:E:53:LEU:HD11	1.82	0.59
1:A:136:SER:HA	2:B:121:PHE:HD2	1.67	0.59
2:E:3:VAL:H	2:E:26:SER:HB3	1.67	0.59
2:E:124:PRO:HB3	2:E:214:PHE:CZ	2.38	0.59
3:C:177:SER:HB2	5:C:323:HOH:O	2.02	0.58
2:B:31:HIS:CE1	2:B:33:ARG:HB2	2.38	0.58
1:D:1:GLN:N	1:D:1:GLN:OE1	2.27	0.58
2:E:121:PHE:C	2:E:122:ILE:HD12	2.24	0.58
2:B:39:LEU:HD23	5:B:465:HOH:O	2.03	0.58
1:D:105:ILE:HD12	2:E:52:LEU:HD21	1.84	0.58
1:A:206:HIS:HB3	1:A:211:THR:OG1	2.04	0.58
2:B:147:ARG:NH2	2:B:168:VAL:HG11	2.18	0.57
1:D:160:TRP:CH2	1:D:202:CYS:HB3	2.37	0.57
1:A:67:ARG:NH2	1:A:87:LYS:HE2	2.19	0.57
1:D:195:LEU:HD21	1:D:219:PRO:HB3	1.86	0.57
1:A:174:ALA:HA	1:A:184:LEU:HB3	1.86	0.57
1:D:190:VAL:HG21	1:D:200:TYR:OH	2.05	0.56
1:A:147:LEU:HD13	5:A:559:HOH:O	2.04	0.56
2:B:124:PRO:HB3	2:B:214:PHE:CE2	2.40	0.56
1:D:165:LEU:HD21	1:D:188:VAL:HG21	1.87	0.56
3:F:138:ASP:OD1	3:F:172:ASN:OD1	2.24	0.56
1:A:1:GLN:OE1	1:A:1:GLN:N	2.31	0.55
1:A:108:TYR:CE1	4:A:401:GOL:H31	2.41	0.55
1:D:160:TRP:CZ3	1:D:202:CYS:HB3	2.41	0.55
2:E:149:ALA:HB2	2:E:203:HIS:HD2	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ASN:ND2	3:C:180:ASN:HB3	2.21	0.55
1:A:105:ILE:HD12	2:B:52:LEU:HD21	1.88	0.55
2:B:200:GLU:OE1	5:B:403:HOH:O	2.18	0.54
2:B:125:PRO:HD3	2:B:137:VAL:HG22	1.88	0.54
2:B:32:SER:HA	2:B:35:ARG:HH21	1.72	0.54
3:C:147:VAL:HG21	3:C:174:LEU:HD13	1.90	0.54
2:B:143:ASN:N	2:B:143:ASN:HD22	2.05	0.54
1:A:91:THR:HA	1:A:115:VAL:O	2.08	0.54
1:A:105:ILE:HD12	2:B:52:LEU:HD22	1.90	0.54
2:B:124:PRO:HB3	2:B:214:PHE:CZ	2.42	0.54
2:B:138:VAL:HG22	2:B:183:THR:HG22	1.90	0.53
3:C:196:ASP:OD2	5:C:301:HOH:O	2.19	0.53
1:D:157:THR:OG1	1:D:205:ASN:HB3	2.08	0.53
1:D:165:LEU:CD2	1:D:188:VAL:HG21	2.38	0.53
2:E:129:GLN:HG2	2:E:134:THR:O	2.08	0.53
2:E:3:VAL:H	2:E:26:SER:CB	2.22	0.53
1:A:125:PRO:HB3	1:A:151:TYR:HB3	1.89	0.53
1:D:91:THR:HA	1:D:115:VAL:O	2.09	0.53
1:D:125:PRO:CB	1:D:151:TYR:HB3	2.37	0.53
2:B:113:ARG:HG2	2:B:114:THR:N	2.24	0.52
2:B:78:THR:O	5:B:404:HOH:O	2.19	0.52
1:D:174:ALA:HA	1:D:184:LEU:HB3	1.92	0.52
1:A:135:LYS:O	5:A:501:HOH:O	2.18	0.52
1:A:136:SER:O	1:A:142:ALA:HA	2.09	0.52
1:A:135:LYS:HB3	5:A:501:HOH:O	2.09	0.52
2:E:18:PRO:HA	2:E:81:ILE:O	2.10	0.52
2:B:206:LEU:HD13	2:B:210:VAL:CG2	2.40	0.52
3:F:116:LYS:C	3:F:116:LYS:HD2	2.31	0.51
1:A:137:THR:HG22	1:A:142:ALA:HB2	1.92	0.51
1:D:152:PHE:CE2	1:D:153:PRO:HB3	2.46	0.51
3:C:132:GLN:O	3:C:136:VAL:HG22	2.10	0.51
1:A:101:PRO:O	1:A:103:VAL:N	2.44	0.51
3:C:140:ALA:O	3:C:144:LYS:HG3	2.11	0.51
1:D:12:LYS:O	1:D:117:VAL:HA	2.11	0.50
1:A:44:GLY:HA3	5:A:537:HOH:O	2.11	0.50
2:B:129:GLN:HG2	2:B:134:THR:O	2.10	0.50
2:B:166:GLU:OE2	2:B:180:LEU:HD11	2.12	0.50
2:B:13:VAL:HG11	2:B:84:VAL:HG21	1.93	0.50
1:D:190:VAL:CG2	1:D:191:PRO:HD2	2.42	0.50
1:A:192:SER:CA	1:A:195:LEU:HD13	2.39	0.49
2:B:145:TYR:CG	2:B:146:PRO:HA	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:172:ASP:HB3	2:E:175:ASP:OD1	2.12	0.49
2:E:61:GLU:OE1	3:F:167:THR:HG21	2.13	0.49
1:D:154:GLU:HB3	1:D:155:PRO:HA	1.95	0.49
1:A:118:SER:HB3	1:A:152:PHE:CZ	2.48	0.49
1:A:201:ILE:HG12	1:A:216:LYS:CB	2.42	0.49
2:B:1:ASP:HB3	5:B:454:HOH:O	2.13	0.49
2:B:71:SER:N	5:B:404:HOH:O	2.28	0.49
1:D:52:ASN:HD21	3:F:180:ASN:HB3	1.76	0.49
2:E:206:LEU:HD13	2:E:210:VAL:HG23	1.95	0.49
2:B:198:ALA:HB2	2:B:213:SER:HB3	1.95	0.49
1:D:207:LYS:N	1:D:208:PRO:HD2	2.27	0.49
1:D:16:ALA:O	1:D:86:LEU:HB2	2.13	0.49
2:E:122:ILE:HG21	2:E:214:PHE:CD2	2.48	0.48
2:E:185:THR:O	2:E:186:LEU:HD23	2.12	0.48
1:D:101:PRO:O	1:D:103:VAL:N	2.43	0.48
1:D:125:PRO:HB3	1:D:151:TYR:CB	2.36	0.48
2:B:39:LEU:HD22	2:B:77:PHE:CD2	2.49	0.48
1:D:20:VAL:CG2	1:D:81:LEU:HB3	2.44	0.48
1:A:125:PRO:CA	1:A:151:TYR:HB3	2.43	0.48
1:A:154:GLU:HB3	1:A:155:PRO:HA	1.96	0.48
2:B:147:ARG:CZ	2:B:168:VAL:HG11	2.44	0.47
2:B:197:TYR:HB2	2:B:214:PHE:CE1	2.49	0.47
2:E:187:SER:O	2:E:191:TYR:N	2.47	0.47
1:D:38:ARG:O	1:D:46:GLU:HG2	2.14	0.47
3:C:111:THR:N	5:C:303:HOH:O	2.47	0.47
1:D:52:ASN:OD1	1:D:55:THR:N	2.40	0.47
1:A:17:SER:HA	1:A:83:ILE:O	2.16	0.46
2:E:120:VAL:HA	2:E:140:LEU:O	2.15	0.46
1:A:82:GLN:HG3	5:A:506:HOH:O	2.16	0.46
1:D:111:GLN:OE1	1:D:111:GLN:N	2.42	0.46
1:D:122:THR:HA	1:D:152:PHE:O	2.14	0.46
1:A:104:VAL:HG12	2:B:97:SER:HB2	1.97	0.46
2:E:145:TYR:CG	2:E:146:PRO:HA	2.51	0.46
2:E:31:HIS:HB2	2:E:38:TYR:HE2	1.79	0.46
2:E:11:LEU:HD21	2:E:19:ALA:HB1	1.96	0.46
3:C:184:ASN:HA	3:C:187:LYS:HE2	1.97	0.46
1:D:130:LEU:HD12	1:D:130:LEU:N	2.31	0.46
1:A:194:SER:HA	1:A:197:THR:HG22	1.98	0.46
2:B:31:HIS:HE1	2:B:33:ARG:HB2	1.77	0.46
3:C:132:GLN:O	3:C:136:VAL:HG13	2.15	0.46
5:A:557:HOH:O	3:C:171:LYS:HB2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:197:TYR:HB2	2:E:214:PHE:CE2	2.50	0.46
2:E:110:GLU:OE1	2:E:178:TYR:OH	2.34	0.46
2:E:125:PRO:HD3	2:E:137:VAL:HG22	1.97	0.46
1:D:39:GLN:O	1:D:92:ALA:HB1	2.16	0.46
3:C:190:CYS:O	3:C:194:ILE:HG13	2.15	0.46
1:D:159:SER:OG	1:D:163:GLY:N	2.48	0.46
3:F:131:LEU:HA	3:F:147:VAL:HG11	1.97	0.46
1:D:190:VAL:HG23	1:D:191:PRO:HD2	1.98	0.45
2:E:122:ILE:HG21	2:E:214:PHE:HD2	1.81	0.45
1:A:122:THR:HA	1:A:152:PHE:O	2.16	0.45
3:C:176:PRO:HB2	5:C:366:HOH:O	2.16	0.45
1:A:108:TYR:HE1	4:A:401:GOL:H31	1.80	0.45
1:A:47:TRP:HZ2	1:A:50:TRP:HD1	1.63	0.45
2:E:198:ALA:CB	2:E:213:SER:HB3	2.44	0.45
2:E:180:LEU:HD23	2:E:181:SER:N	2.32	0.45
2:E:124:PRO:HB3	2:E:214:PHE:CE1	2.51	0.45
2:B:67:ARG:NH1	2:B:88:ASP:OD2	2.49	0.45
1:D:195:LEU:HD21	1:D:219:PRO:CB	2.46	0.45
1:A:111:GLN:OE1	1:A:111:GLN:N	2.41	0.45
1:A:175:VAL:HG12	1:A:183:SER:O	2.16	0.45
2:B:56:TRP:O	2:B:57:ALA:HB3	2.17	0.45
1:D:35:HIS:ND1	1:D:50:TRP:HB3	2.32	0.45
2:E:56:TRP:O	2:E:57:ALA:HB3	2.17	0.45
1:A:149:LYS:HE3	1:A:177:GLN:HE22	1.81	0.44
1:A:118:SER:HB3	1:A:152:PHE:HZ	1.81	0.44
1:D:17:SER:HA	1:D:83:ILE:O	2.18	0.44
1:A:125:PRO:HA	1:A:151:TYR:HB3	2.00	0.44
2:B:2:ILE:HA	2:B:26:SER:OG	2.18	0.44
2:E:122:ILE:HD11	2:E:199:CYS:SG	2.57	0.44
2:E:95:LYS:HB2	2:E:103:PHE:CD2	2.52	0.44
2:E:43:LEU:O	2:E:51:GLN:HG2	2.18	0.44
3:F:148:LYS:O	3:F:152:GLU:HG3	2.18	0.44
2:B:185:THR:O	2:B:186:LEU:HD23	2.17	0.43
1:A:123:LYS:HE2	1:A:181:LEU:CD1	2.48	0.43
1:A:89:GLU:CD	2:E:66:ASP:HB2	2.38	0.43
2:E:13:VAL:HG11	2:E:84:VAL:CG2	2.47	0.43
1:A:159:SER:OG	1:A:163:GLY:N	2.51	0.43
2:B:165:GLN:O	2:B:182:SER:HA	2.18	0.43
2:E:67:ARG:O	2:E:81:ILE:HA	2.18	0.43
1:D:89:GLU:N	1:D:89:GLU:OE1	2.51	0.43
2:E:99:ASN:O	2:E:100:LEU:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:17:GLU:OE1	2:E:17:GLU:HA	2.19	0.43
1:A:90:ASP:HA	5:A:502:HOH:O	2.19	0.43
3:C:131:LEU:HD11	3:C:169:VAL:CG2	2.49	0.43
1:D:35:HIS:CG	1:D:50:TRP:HB3	2.54	0.43
2:E:172:ASP:OD2	2:E:174:LYS:HB2	2.18	0.43
2:E:122:ILE:HD13	2:E:212:LYS:CB	2.49	0.43
2:B:23:CYS:HB2	2:B:41:TRP:CH2	2.54	0.43
1:D:151:TYR:OH	1:D:184:LEU:HD23	2.18	0.43
2:E:141:LEU:HD12	2:E:141:LEU:N	2.34	0.43
1:A:38:ARG:O	1:A:46:GLU:HG2	2.19	0.42
2:B:130:LEU:O	2:B:188:LYS:HD2	2.19	0.42
2:B:67:ARG:HH12	2:B:88:ASP:CG	2.21	0.42
3:C:131:LEU:HD11	3:C:169:VAL:HG21	2.00	0.42
3:F:190:CYS:O	3:F:194:ILE:HG13	2.19	0.42
2:B:163:ASN:O	2:B:184:LEU:HD12	2.20	0.42
2:B:128:GLU:N	2:B:128:GLU:OE2	2.44	0.42
2:B:96:GLN:O	2:B:101:TYR:HA	2.19	0.42
1:A:38:ARG:NH1	5:A:502:HOH:O	2.19	0.42
1:A:184:LEU:HA	5:A:563:HOH:O	2.19	0.42
3:F:165:LEU:O	3:F:169:VAL:HG23	2.20	0.42
1:D:91:THR:CG2	1:D:116:THR:HA	2.36	0.42
1:A:125:PRO:CB	1:A:151:TYR:HB3	2.50	0.42
3:F:133:GLN:O	3:F:136:VAL:HG22	2.20	0.42
2:B:95:LYS:HB3	2:B:95:LYS:HE2	1.90	0.41
3:F:147:VAL:HG21	3:F:174:LEU:HD13	2.01	0.41
1:D:15:GLY:N	1:D:86:LEU:O	2.53	0.41
3:C:131:LEU:HA	3:C:147:VAL:HG11	2.02	0.41
3:F:116:LYS:O	3:F:119:ILE:HG22	2.21	0.41
1:D:206:HIS:NE2	1:D:208:PRO:HG2	2.34	0.41
2:B:187:SER:O	2:B:191:TYR:N	2.52	0.41
3:C:173:ASN:HA	5:C:321:HOH:O	2.20	0.41
2:E:24:LYS:HA	2:E:75:THR:O	2.21	0.41
1:A:156:VAL:CG1	1:A:184:LEU:HD21	2.51	0.41
2:B:79:LEU:C	2:B:79:LEU:HD13	2.41	0.41
2:E:180:LEU:HD23	2:E:180:LEU:C	2.40	0.41
2:B:113:ARG:CG	2:B:114:THR:N	2.84	0.41
2:E:39:LEU:HD12	2:E:95:LYS:O	2.20	0.41
3:F:193:LYS:HA	3:F:193:LYS:HD3	1.84	0.41
2:B:99:ASN:O	2:B:100:LEU:HB2	2.20	0.41
1:D:129:PRO:C	1:D:130:LEU:HD12	2.42	0.41
1:D:4:LEU:HD23	1:D:24:ALA:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:SER:N	2:E:121:PHE:HD1	2.18	0.40
1:A:184:LEU:C	1:A:184:LEU:HD12	2.42	0.40
2:E:13:VAL:HG21	2:E:19:ALA:HB2	2.04	0.40
1:A:65:LYS:HB2	1:A:65:LYS:HE3	1.87	0.40
2:E:38:TYR:CZ	3:F:186:PHE:HD1	2.40	0.40

All (3) 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)
5:C:360:HOH:O	5:F:308:HOH:O[3_454]	1.96	0.24
5:A:550:HOH:O	5:C:303:HOH:O[3_544]	2.08	0.12
5:C:333:HOH:O	5:F:343:HOH:O[3_554]	2.10	0.10

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	219/232 (94%)	212 (97%)	7 (3%)	0	100	100
1	D	214/232 (92%)	209 (98%)	5 (2%)	0	100	100
2	B	217/219 (99%)	210 (97%)	7 (3%)	0	100	100
2	E	216/219 (99%)	209 (97%)	7 (3%)	0	100	100
3	C	90/98 (92%)	87 (97%)	3 (3%)	0	100	100
3	F	90/98 (92%)	87 (97%)	3 (3%)	0	100	100
All	All	1046/1098 (95%)	1014 (97%)	32 (3%)	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	185/197 (94%)	185 (100%)	0	100	100
1	D	176/197 (89%)	176 (100%)	0	100	100
2	B	188/194 (97%)	186 (99%)	2 (1%)	73	87
2	E	186/194 (96%)	185 (100%)	1 (0%)	88	95
3	C	81/87 (93%)	80 (99%)	1 (1%)	71	85
3	F	80/87 (92%)	79 (99%)	1 (1%)	69	84
All	All	896/956 (94%)	891 (99%)	5 (1%)	86	94

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	9	LEU
2	B	143	ASN
3	C	132	GLN
2	E	60	ARG
3	F	116	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	GLN
2	B	143	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

3 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	$\# Z > 2$
4	GOL	B	301	-	5,5,5	0.72	0	5,5,5	0.88	0
4	GOL	E	301	-	5,5,5	0.78	0	5,5,5	0.95	0
4	GOL	A	401	-	5,5,5	0.71	0	5,5,5	0.85	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
4	GOL	B	301	-	-	4/4/4/4	-
4	GOL	E	301	-	-	2/4/4/4	-
4	GOL	A	401	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	GOL	O1-C1-C2-C3
4	B	301	GOL	C1-C2-C3-O3
4	E	301	GOL	O1-C1-C2-C3
4	A	401	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
4	B	301	GOL	O2-C2-C3-O3
4	A	401	GOL	O2-C2-C3-O3
4	B	301	GOL	O1-C1-C2-O2
4	E	301	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	GOL	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 ⓘ

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	221/232 (95%)	-0.04	2 (0%) 84 82	28, 62, 100, 143	0
1	D	218/232 (93%)	0.77	34 (15%) 2 1	42, 110, 152, 169	0
2	B	219/219 (100%)	-0.22	0 100 100	34, 63, 112, 133	0
2	E	218/219 (99%)	0.42	20 (9%) 9 8	46, 96, 169, 251	0
3	C	92/98 (93%)	-0.29	0 100 100	29, 47, 86, 100	0
3	F	92/98 (93%)	-0.06	1 (1%) 80 79	34, 54, 84, 108	0
All	All	1060/1098 (96%)	0.16	57 (5%) 25 24	28, 72, 145, 251	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	137	VAL	7.4
1	D	217	VAL	6.3
2	E	133	GLY	5.5
2	E	186	LEU	4.8
1	A	221	SER	4.4
1	D	199	THR	4.3
2	E	185	THR	4.2
2	E	136	SER	4.1
1	D	68	PHE	4.1
2	E	196	VAL	3.9
1	D	204	VAL	3.9
1	D	182	TYR	3.7
2	E	134	THR	3.6
1	D	81	LEU	3.6
2	E	189	ALA	3.6
2	E	153	TRP	3.5
1	D	86	LEU	3.4
2	E	159	LEU	3.0
1	D	147	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	85	SER	2.9
2	E	210	VAL	2.9
1	D	84	SER	2.9
1	D	72	LEU	2.8
1	D	166	THR	2.8
2	E	135	ALA	2.8
1	D	118	SER	2.8
1	D	115	VAL	2.8
1	D	49	GLY	2.8
1	D	88	ALA	2.7
1	A	137	THR	2.6
2	E	187	SER	2.6
2	E	216	ARG	2.5
2	E	158	ALA	2.5
1	D	175	VAL	2.5
2	E	132	SER	2.5
1	D	144	LEU	2.5
1	D	156	VAL	2.5
1	D	218	GLU	2.4
1	D	20	VAL	2.4
1	D	82	GLN	2.4
1	D	138	SER	2.4
3	F	138	ASP	2.4
2	E	190	ASP	2.4
1	D	190	VAL	2.3
1	D	186	SER	2.3
2	E	130	LEU	2.2
1	D	127	VAL	2.2
1	D	119	SER	2.2
1	D	53	THR	2.1
1	D	160	TRP	2.1
1	D	77	SER	2.1
1	D	126	SER	2.0
2	E	140	LEU	2.0
1	D	213	VAL	2.0
1	D	128	PHE	2.0
2	E	127	ASP	2.0
1	D	153	PRO	2.0

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

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. 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(\AA^2)	Q<0.9
4	GOL	B	301	6/6	0.69	0.24	83,100,117,117	0
4	GOL	E	301	6/6	0.82	0.34	118,141,150,150	0
4	GOL	A	401	6/6	0.90	0.23	97,117,125,125	0

6.5 Other polymers [i](#)

There are no such residues in this entry.