Full wwPDB X-ray Structure Validation Report (i)

Sep 12, 2023 – 11:46 PM EDT

PDB ID	:	4NEE
Title	:	crystal structure of AP-2 alpha/simga2 complex bound to HIV-1 Nef
Authors	:	Hurley, J.H.; Bonifacino, J.S.; Ren, X.; Park, S.Y.
Deposited on	:	2013-10-29
Resolution	:	2.88 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 (i) 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 (1)) were used in the production of this report:

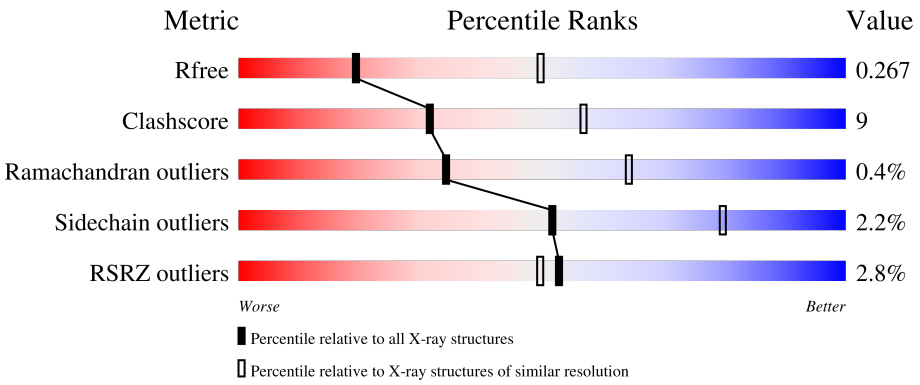
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

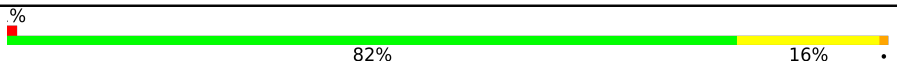
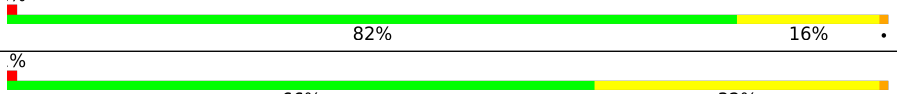
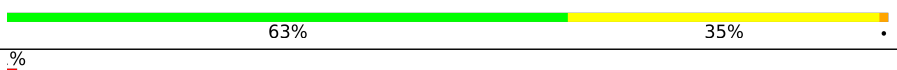
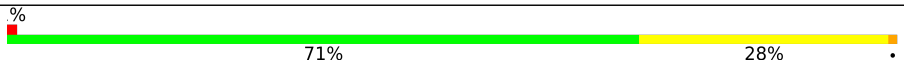
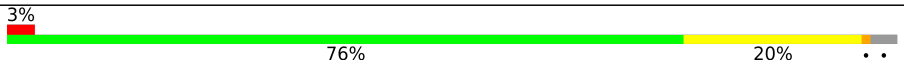
The reported resolution of this entry is 2.88 Å.

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	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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 <=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	D	142	
1	F	142	
1	I	142	
1	L	142	
2	A	398	

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Mol	Chain	Length	Quality of chain
2	B	398	2% 78% 19%
2	G	398	2% 79% 17%
2	J	398	2% 76% 21%
3	C	155	5% 66% 23% 10%
3	E	155	% 66% 23% 10%
3	H	155	14% 66% 25% 8%
3	K	155	6% 73% 19% 7%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 21530 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 AP-2 complex subunit sigma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	142	1199	778	200	214	7	0	0	0
1	L	142	1199	778	200	214	7	0	0	0
1	F	142	1199	778	200	214	7	0	0	0
1	I	142	1199	778	200	214	7	0	0	0

- Molecule 2 is a protein called AP-2 complex subunit alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	G	388	3047	1931	536	564	16	0	0	0
2	B	388	3047	1931	536	564	16	0	0	0
2	J	388	3047	1931	536	564	16	0	0	0
2	A	385	3025	1918	532	559	16	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	-1	GLY	-	cloning artifact	UNP Q66HM2
G	0	ALA	-	cloning artifact	UNP Q66HM2
B	-1	GLY	-	cloning artifact	UNP Q66HM2
B	0	ALA	-	cloning artifact	UNP Q66HM2
J	-1	GLY	-	cloning artifact	UNP Q66HM2
J	0	ALA	-	cloning artifact	UNP Q66HM2
A	-1	GLY	-	cloning artifact	UNP Q66HM2
A	0	ALA	-	cloning artifact	UNP Q66HM2

- Molecule 3 is a protein called Protein Nef.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	E	139	1134	732	197	202	3	0	0	0
			Total	C	N	O	S			
3	H	142	1143	737	199	204	3	0	0	0
			Total	C	N	O	S			
3	K	144	1157	745	202	207	3	0	0	0
			Total	C	N	O	S			
3	C	139	1134	732	197	202	3	0	0	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	48	GLY	-	expression tag	UNP Q90VU7
E	49	UNK	-	expression tag	UNP Q90VU7
E	50	UNK	-	expression tag	UNP Q90VU7
E	51	UNK	-	expression tag	UNP Q90VU7
E	52	UNK	-	expression tag	UNP Q90VU7
E	53	UNK	-	expression tag	UNP Q90VU7
E	54	UNK	-	expression tag	UNP Q90VU7
E	55	UNK	-	expression tag	UNP Q90VU7
E	56	UNK	-	expression tag	UNP Q90VU7
E	57	UNK	-	expression tag	UNP Q90VU7
E	58	UNK	-	expression tag	UNP Q90VU7
E	59	UNK	-	expression tag	UNP Q90VU7
E	60	UNK	-	expression tag	UNP Q90VU7
E	61	UNK	-	expression tag	UNP Q90VU7
H	54	GLY	-	expression tag	UNP Q90VU7
H	55	UNK	-	expression tag	UNP Q90VU7
H	56	UNK	-	expression tag	UNP Q90VU7
H	57	UNK	-	expression tag	UNP Q90VU7
H	58	UNK	-	expression tag	UNP Q90VU7
H	59	UNK	-	expression tag	UNP Q90VU7
H	60	UNK	-	expression tag	UNP Q90VU7
H	61	UNK	-	expression tag	UNP Q90VU7
H	62	UNK	-	expression tag	UNP Q90VU7
H	63	UNK	-	expression tag	UNP Q90VU7
H	64	UNK	-	expression tag	UNP Q90VU7
H	65	UNK	-	expression tag	UNP Q90VU7
H	66	UNK	-	expression tag	UNP Q90VU7
H	67	UNK	-	expression tag	UNP Q90VU7
K	54	GLY	-	expression tag	UNP Q90VU7
K	55	UNK	-	expression tag	UNP Q90VU7

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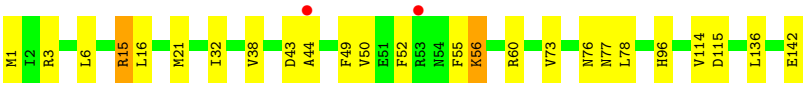
Chain	Residue	Modelled	Actual	Comment	Reference
K	56	UNK	-	expression tag	UNP Q90VU7
K	57	UNK	-	expression tag	UNP Q90VU7
K	58	UNK	-	expression tag	UNP Q90VU7
K	59	UNK	-	expression tag	UNP Q90VU7
K	60	UNK	-	expression tag	UNP Q90VU7
K	61	UNK	-	expression tag	UNP Q90VU7
K	62	UNK	-	expression tag	UNP Q90VU7
K	63	UNK	-	expression tag	UNP Q90VU7
K	64	UNK	-	expression tag	UNP Q90VU7
K	65	UNK	-	expression tag	UNP Q90VU7
K	66	UNK	-	expression tag	UNP Q90VU7
K	67	UNK	-	expression tag	UNP Q90VU7
C	48	GLY	-	expression tag	UNP Q90VU7
C	49	UNK	-	expression tag	UNP Q90VU7
C	50	UNK	-	expression tag	UNP Q90VU7
C	51	UNK	-	expression tag	UNP Q90VU7
C	52	UNK	-	expression tag	UNP Q90VU7
C	53	UNK	-	expression tag	UNP Q90VU7
C	54	UNK	-	expression tag	UNP Q90VU7
C	55	UNK	-	expression tag	UNP Q90VU7
C	56	UNK	-	expression tag	UNP Q90VU7
C	57	UNK	-	expression tag	UNP Q90VU7
C	58	UNK	-	expression tag	UNP Q90VU7
C	59	UNK	-	expression tag	UNP Q90VU7
C	60	UNK	-	expression tag	UNP Q90VU7
C	61	UNK	-	expression tag	UNP Q90VU7

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.

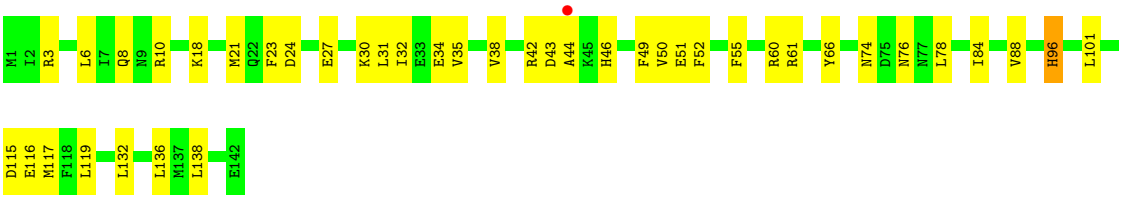
• Molecule 1: AP-2 complex subunit sigma

Chain D:



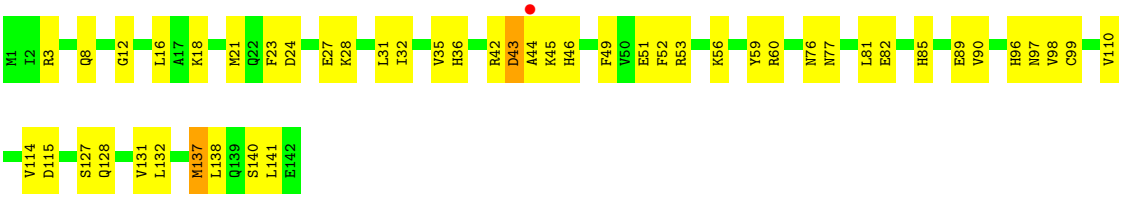
• Molecule 1: AP-2 complex subunit sigma

Chain L:



• Molecule 1: AP-2 complex subunit sigma

Chain F:

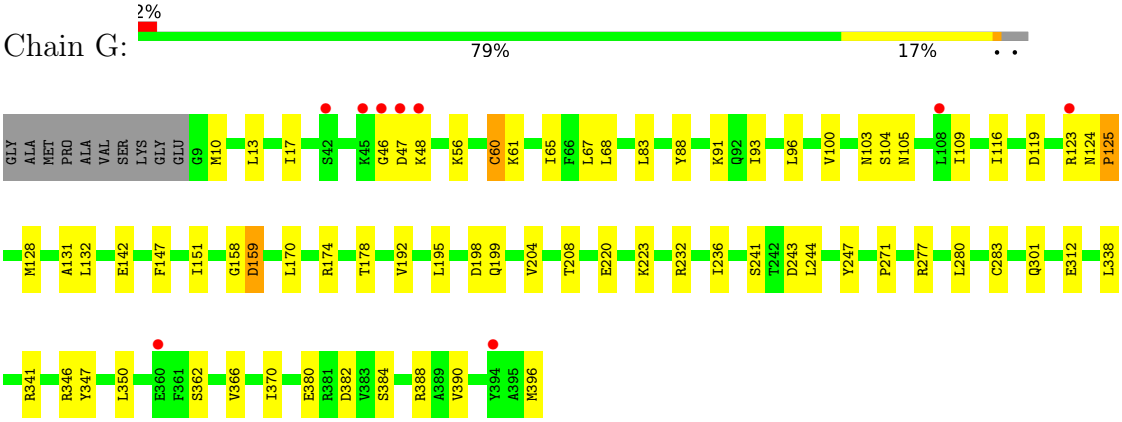


• Molecule 1: AP-2 complex subunit sigma

Chain I:

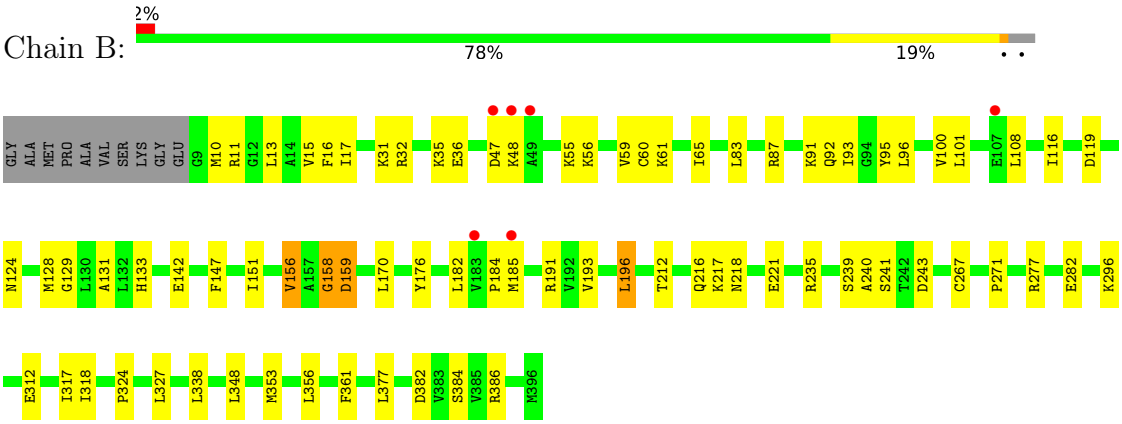
• Molecule 2: AP-2 complex subunit alpha-2

Chain G:



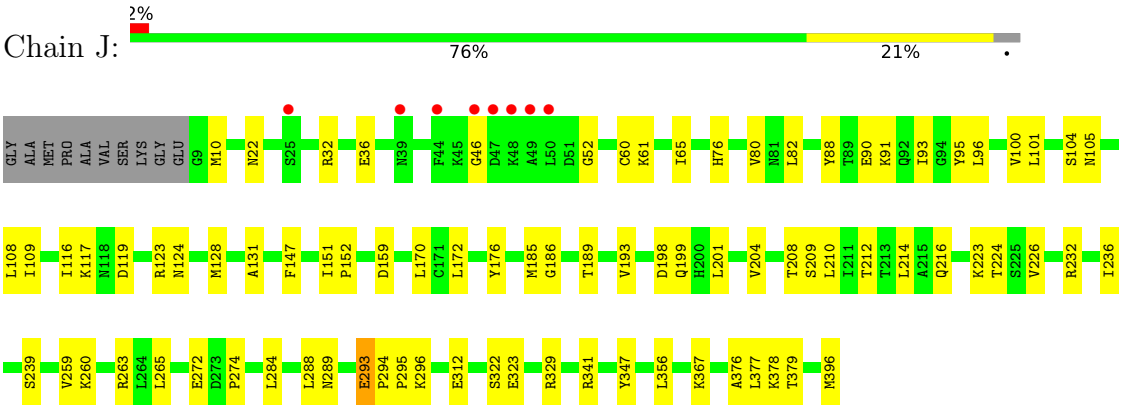
• Molecule 2: AP-2 complex subunit alpha-2

Chain B:



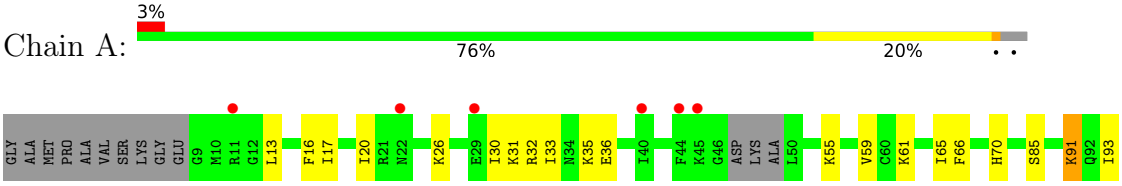
• Molecule 2: AP-2 complex subunit alpha-2

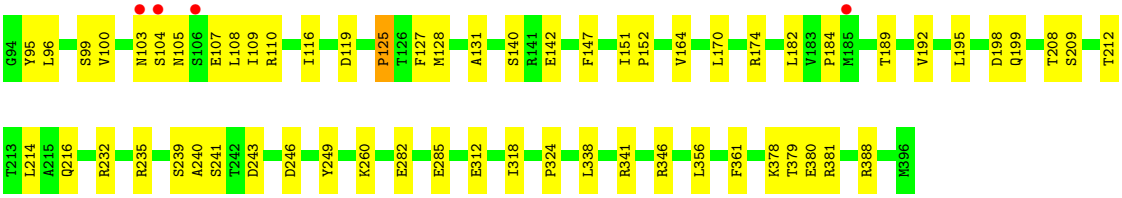
Chain J:



• Molecule 2: AP-2 complex subunit alpha-2

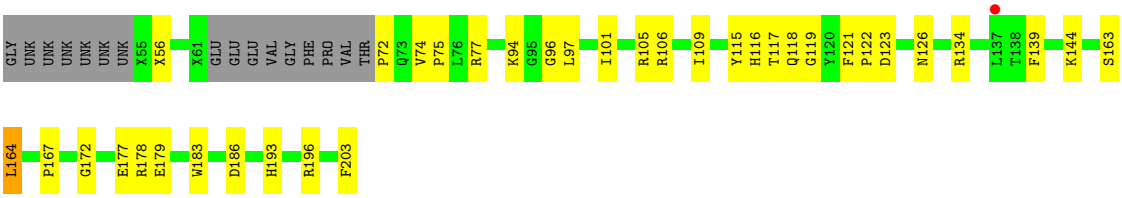
Chain A:





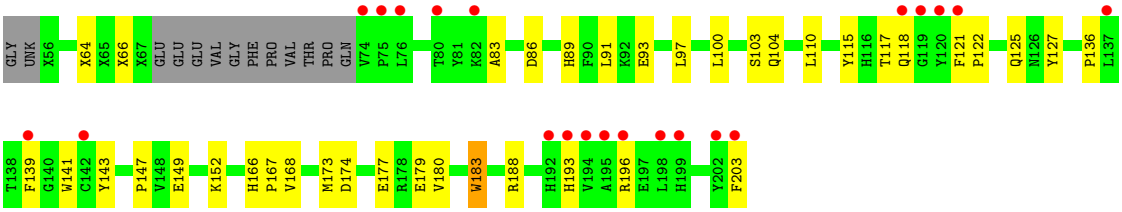
• Molecule 3: Protein Nef

Chain E:



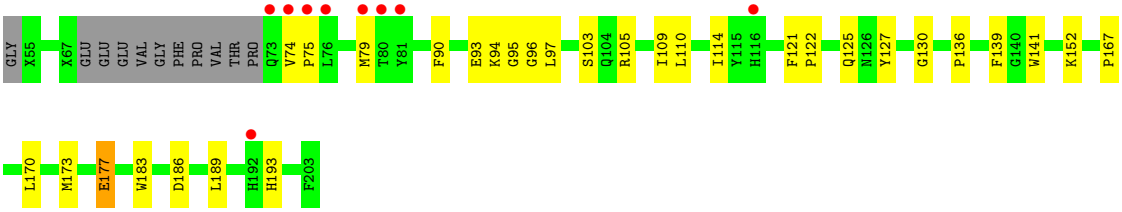
• Molecule 3: Protein Nef

Chain H:



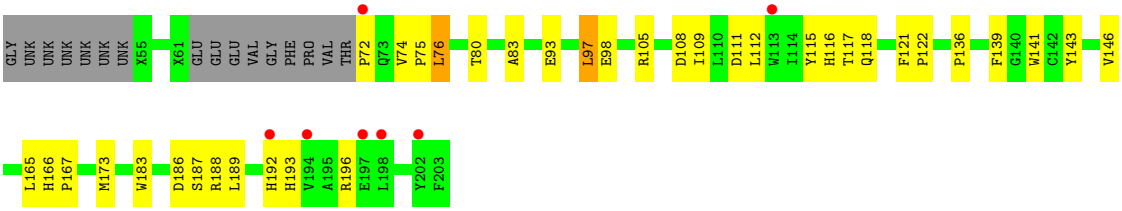
• Molecule 3: Protein Nef

Chain K:



• Molecule 3: Protein Nef

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	109.56Å 168.03Å 200.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.88 – 2.88 48.88 – 2.88	Depositor EDS
% Data completeness (in resolution range)	95.5 (48.88-2.88) 94.6 (48.88-2.88)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.97 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, R_{free}	0.219 , 0.267 0.222 , 0.267	Depositor DCC
R_{free} test set	4014 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	35.7	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 39.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.42$, $< L^2 > = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	21530	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.36% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	D	0.81	0/1223	0.60	0/1650
1	F	0.73	0/1223	0.60	0/1650
1	I	0.87	0/1223	0.64	0/1650
1	L	0.80	0/1223	0.65	0/1650
2	A	0.61	0/3076	0.57	0/4161
2	B	0.69	1/3099 (0.0%)	0.59	1/4193 (0.0%)
2	G	0.70	0/3099	0.57	0/4193
2	J	0.68	1/3099 (0.0%)	0.59	0/4193
3	C	0.71	0/1139	0.59	0/1551
3	E	0.70	0/1139	0.59	0/1551
3	H	0.62	0/1122	0.58	0/1528
3	K	0.64	0/1131	0.55	0/1540
All	All	0.70	2/21796 (0.0%)	0.59	1/29510 (0.0%)

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	F	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	60	CYS	CB-SG	-5.15	1.73	1.81
2	B	267	CYS	CB-SG	-5.05	1.73	1.81

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	158	GLY	N-CA-C	-5.77	98.67	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	43	ASP	Peptide

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	D	1199	0	1195	18	0
1	F	1199	0	1195	37	0
1	I	1199	0	1195	36	0
1	L	1199	0	1195	31	0
2	A	3025	0	3105	52	0
2	B	3047	0	3128	52	0
2	G	3047	0	3128	46	0
2	J	3047	0	3128	59	0
3	C	1134	0	1058	25	0
3	E	1134	0	1058	26	0
3	H	1143	0	1047	30	0
3	K	1157	0	1056	19	0
All	All	21530	0	21488	386	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 9.

All (386) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:209:SER:OG	2:A:260:LYS:NZ	2.04	0.91
1:F:56:LYS:NZ	1:F:77:ASN:OD1	2.09	0.86
1:L:76:ASN:HD22	2:J:312:GLU:HG2	1.40	0.86
2:A:95:TYR:OH	2:A:119:ASP:OD2	1.95	0.83
3:E:56:UNK:HA	3:E:94:LYS:HE2	1.61	0.83
1:F:46:HIS:O	3:E:178:ARG:NH1	2.13	0.82
3:C:193:HIS:HD2	3:C:196:ARG:H	1.28	0.81
1:D:60:ARG:NH2	3:C:173:MET:SD	2.55	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:43:ASP:HB3	1:F:45:LYS:N	1.98	0.79
3:E:134:ARG:NH2	3:E:179:GLU:OE2	2.17	0.78
3:H:193:HIS:HD2	3:H:196:ARG:H	1.30	0.78
1:I:56:LYS:NZ	1:I:77:ASN:OD1	2.16	0.77
1:L:18:LYS:NZ	1:L:115:ASP:OD2	2.15	0.77
2:J:209:SER:OG	2:J:260:LYS:NZ	2.18	0.76
1:D:38:VAL:HG11	1:D:50:VAL:HG11	1.71	0.73
2:G:56:LYS:HG3	2:G:93:ILE:HG13	1.70	0.73
2:B:156:VAL:HG22	2:B:191:ARG:HB2	1.71	0.73
1:D:1:MET:SD	1:D:3:ARG:NH1	2.61	0.73
2:B:382:ASP:OD2	2:B:384:SER:OG	2.07	0.72
3:H:149:GLU:HB2	3:H:152:LYS:HG3	1.72	0.72
2:G:301:GLN:HE21	3:H:173:MET:HE1	1.53	0.71
2:B:91:LYS:HE2	2:B:119:ASP:OD1	1.89	0.71
2:A:125:PRO:O	2:A:127:PHE:N	2.21	0.71
3:H:64:UNK:O	3:H:66:UNK:HA	1.91	0.71
2:A:318:ILE:HG12	2:A:356:LEU:HD13	1.75	0.69
1:L:38:VAL:HG11	1:L:50:VAL:HG11	1.74	0.69
3:C:80:THR:H	3:C:83:ALA:HB3	1.58	0.69
2:B:240:ALA:HB1	2:A:282:GLU:OE1	1.94	0.68
2:J:95:TYR:OH	2:J:119:ASP:OD1	2.04	0.67
2:A:235:ARG:O	2:A:239:SER:OG	2.12	0.67
1:D:16:LEU:HD21	1:D:114:VAL:HG21	1.77	0.67
1:D:43:ASP:OD1	1:D:44:ALA:N	2.26	0.67
2:J:116:ILE:HG23	2:J:131:ALA:HB1	1.76	0.66
2:A:85:SER:O	2:A:91:LYS:NZ	2.21	0.66
2:A:208:THR:O	2:A:212:THR:OG1	2.13	0.66
1:I:38:VAL:HG11	1:I:50:VAL:HG11	1.78	0.65
3:C:72:PRO:HD2	3:C:116:HIS:HA	1.78	0.65
2:G:91:LYS:HE2	2:G:119:ASP:OD1	1.96	0.65
2:A:174:ARG:O	2:A:174:ARG:NH1	2.30	0.65
2:G:116:ILE:HG23	2:G:131:ALA:HB1	1.77	0.64
1:I:1:MET:SD	1:I:3:ARG:NH2	2.71	0.64
1:L:76:ASN:ND2	2:J:312:GLU:HG2	2.11	0.64
2:A:198:ASP:OD1	2:A:199:GLN:N	2.31	0.64
3:C:105:ARG:O	3:C:109:ILE:HG13	1.98	0.64
2:J:236:ILE:O	2:J:239:SER:OG	2.10	0.63
2:B:377:LEU:O	2:B:386:ARG:NH2	2.31	0.63
1:D:56:LYS:HE2	1:D:73:VAL:HA	1.81	0.63
3:K:125:GLN:NE2	3:K:127:TYR:OH	2.26	0.63
1:F:21:MET:HE2	1:F:23:PHE:CZ	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:95:TYR:OH	2:B:119:ASP:OD2	2.16	0.62
1:F:44:ALA:O	3:E:178:ARG:NH1	2.33	0.62
3:E:101:ILE:O	3:E:106:ARG:NH1	2.33	0.61
1:I:30:LYS:O	1:I:34:GLU:HG3	2.01	0.61
3:H:125:GLN:NE2	3:H:127:TYR:OH	2.25	0.61
1:D:76:ASN:ND2	2:B:312:GLU:OE1	2.31	0.60
2:A:346:ARG:NH1	2:A:380:GLU:OE1	2.35	0.60
2:A:192:VAL:O	2:A:195:LEU:HB2	2.01	0.60
2:A:240:ALA:HB3	2:A:243:ASP:OD2	2.02	0.59
1:F:3:ARG:HD2	1:F:21:MET:SD	2.43	0.59
2:J:61:LYS:O	2:J:65:ILE:HG13	2.03	0.58
1:I:3:ARG:HD2	1:I:21:MET:SD	2.42	0.58
1:I:127:SER:O	1:I:131:VAL:HG23	2.03	0.58
2:G:346:ARG:NH1	2:G:380:GLU:OE1	2.37	0.58
3:K:103:SER:HB3	3:K:167:PRO:HB3	1.85	0.58
1:L:101:LEU:HD21	3:C:187:SER:OG	2.04	0.58
2:B:235:ARG:O	2:B:239:SER:OG	2.21	0.58
2:A:105:ASN:O	2:A:109:ILE:HG13	2.04	0.57
2:J:198:ASP:OD1	2:J:199:GLN:N	2.37	0.57
1:D:6:LEU:HD11	1:D:32:ILE:HD13	1.86	0.57
1:F:76:ASN:ND2	2:A:312:GLU:OE1	2.26	0.57
2:G:382:ASP:OD2	2:G:384:SER:OG	2.19	0.57
1:F:98:VAL:HG13	3:E:164:LEU:HD22	1.85	0.57
2:B:185:MET:HG3	2:B:218:ASN:HD22	1.69	0.57
3:C:193:HIS:CD2	3:C:196:ARG:H	2.18	0.57
1:I:46:HIS:ND1	2:G:382:ASP:OD1	2.38	0.57
2:G:220:GLU:O	2:G:223:LYS:HG3	2.05	0.57
2:A:147:PHE:O	2:A:151:ILE:HG12	2.05	0.57
3:H:166:HIS:HD2	3:H:168:VAL:HB	1.70	0.57
1:I:24:ASP:OD1	1:I:27:GLU:N	2.33	0.57
1:I:86:ASN:OD1	1:I:128:GLN:NE2	2.28	0.56
2:G:390:VAL:HG21	2:J:378:LYS:HB2	1.87	0.56
2:B:240:ALA:HB3	2:B:243:ASP:OD1	2.04	0.56
1:L:42:ARG:NH2	1:L:46:HIS:HB3	2.20	0.56
3:K:130:GLY:HA3	3:K:177:GLU:HG3	1.86	0.56
2:B:142:GLU:OE1	2:B:142:GLU:N	2.38	0.56
1:F:137:MET:O	1:F:141:LEU:HD13	2.05	0.56
1:F:18:LYS:NZ	1:F:115:ASP:OD1	2.39	0.56
1:I:42:ARG:NH1	1:I:46:HIS:HB3	2.21	0.56
1:L:34:GLU:O	1:L:38:VAL:HG23	2.05	0.55
2:J:198:ASP:O	2:J:232:ARG:NH2	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:6:LEU:HD11	1:L:32:ILE:HD13	1.88	0.55
1:L:21:MET:HB3	1:L:23:PHE:CE2	2.41	0.55
1:L:138:LEU:HD22	2:J:88:TYR:CZ	2.41	0.55
1:F:45:LYS:HD3	2:A:381:ARG:HG2	1.88	0.55
1:F:99:CYS:HA	3:E:164:LEU:HD13	1.89	0.55
2:J:259:VAL:HG13	2:J:312:GLU:HG3	1.88	0.55
1:L:3:ARG:HD2	1:L:21:MET:SD	2.46	0.54
2:G:142:GLU:OE1	2:G:142:GLU:N	2.34	0.54
2:B:56:LYS:HG3	2:B:93:ILE:HG13	1.89	0.54
3:K:136:PRO:HB3	3:K:141:TRP:HD1	1.71	0.54
2:B:241:SER:HB3	2:A:282:GLU:OE2	2.07	0.54
2:J:289:ASN:OD1	2:J:329:ARG:NH2	2.39	0.54
2:A:32:ARG:HH12	2:A:36:GLU:HB2	1.71	0.54
1:L:61:ARG:HD2	1:L:66:TYR:CE1	2.42	0.54
2:G:271:PRO:HD2	2:G:277:ARG:HG3	1.89	0.54
3:H:139:PHE:HB2	3:H:193:HIS:HD1	1.73	0.54
1:F:16:LEU:HD21	1:F:114:VAL:HG21	1.89	0.54
2:A:142:GLU:OE1	2:A:142:GLU:N	2.36	0.54
1:D:78:LEU:HD13	2:B:348:LEU:HD11	1.90	0.54
1:F:42:ARG:CZ	1:F:46:HIS:HB3	2.38	0.54
3:K:121:PHE:CD1	3:K:122:PRO:HD2	2.42	0.54
2:A:61:LYS:O	2:A:65:ILE:HG13	2.08	0.53
3:C:186:ASP:HB3	3:C:189:LEU:HG	1.90	0.53
1:D:56:LYS:HD2	1:D:77:ASN:OD1	2.08	0.53
1:I:56:LYS:HG2	1:I:73:VAL:HA	1.90	0.53
2:J:100:VAL:HB	2:J:101:LEU:HD12	1.90	0.53
3:H:193:HIS:CD2	3:H:196:ARG:H	2.18	0.53
3:K:93:GLU:C	3:K:95:GLY:H	2.11	0.53
2:J:117:LYS:HG3	2:J:147:PHE:HE1	1.73	0.52
2:B:282:GLU:OE2	2:A:240:ALA:HB1	2.09	0.52
2:G:60:CYS:SG	2:G:93:ILE:HD12	2.50	0.52
2:G:147:PHE:O	2:G:151:ILE:HG12	2.10	0.52
1:F:24:ASP:N	1:F:27:GLU:OE1	2.34	0.52
1:F:97:ASN:ND2	3:E:163:SER:OG	2.43	0.52
2:G:198:ASP:OD1	2:G:199:GLN:N	2.43	0.52
2:G:198:ASP:O	2:G:232:ARG:NH2	2.43	0.52
2:G:13:LEU:O	2:G:17:ILE:HG13	2.10	0.52
1:D:142:GLU:OE1	2:B:87:ARG:NH1	2.43	0.52
1:L:60:ARG:NH2	3:K:173:MET:SD	2.63	0.52
3:K:139:PHE:HB2	3:K:193:HIS:CD2	2.45	0.52
2:B:124:ASN:O	2:B:128:MET:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:204:VAL:O	2:G:208:THR:OG1	2.23	0.51
3:E:72:PRO:HG3	3:E:116:HIS:CE1	2.46	0.51
2:A:16:PHE:O	2:A:20:ILE:HG12	2.11	0.51
3:E:139:PHE:HB2	3:E:193:HIS:CD2	2.46	0.51
1:F:43:ASP:HB3	1:F:44:ALA:C	2.31	0.51
2:G:96:LEU:O	2:G:100:VAL:HG23	2.10	0.51
2:J:22:ASN:OD1	3:C:188:ARG:HD3	2.10	0.51
3:K:186:ASP:HB3	3:K:189:LEU:HG	1.92	0.51
1:F:60:ARG:HH21	1:F:81:LEU:HG	1.75	0.51
1:F:43:ASP:CB	1:F:44:ALA:HB3	2.40	0.51
2:J:76:HIS:O	2:J:80:VAL:HG23	2.11	0.51
2:J:341:ARG:NH1	3:K:177:GLU:OE2	2.41	0.51
3:K:105:ARG:O	3:K:109:ILE:HG13	2.11	0.51
3:H:91:LEU:HD12	3:H:141:TRP:HH2	1.74	0.50
3:C:117:THR:OG1	3:C:118:GLN:HG3	2.12	0.50
2:A:378:LYS:HG3	2:A:379:THR:HG23	1.93	0.50
3:E:105:ARG:O	3:E:109:ILE:HG13	2.11	0.50
1:I:61:ARG:HD3	1:I:66:TYR:CE1	2.46	0.50
2:G:124:ASN:O	2:G:128:MET:HG3	2.12	0.50
2:G:350:LEU:HB3	2:G:388:ARG:O	2.12	0.50
3:C:139:PHE:HB2	3:C:193:HIS:ND1	2.26	0.50
2:J:80:VAL:HG21	2:J:108:LEU:HD22	1.93	0.50
1:L:76:ASN:H	2:J:263:ARG:HH22	1.58	0.50
1:F:42:ARG:HD3	1:F:59:TYR:OH	2.11	0.50
1:D:3:ARG:HG3	1:D:55:PHE:CZ	2.47	0.49
2:B:318:ILE:HG12	2:B:356:LEU:HD13	1.95	0.49
3:C:115:TYR:CD1	3:C:122:PRO:HD3	2.47	0.49
2:A:31:LYS:O	2:A:35:LYS:HG3	2.13	0.49
1:D:115:ASP:HB3	2:B:133:HIS:CD2	2.47	0.49
1:I:138:LEU:HD23	1:I:141:LEU:HD12	1.94	0.49
3:E:123:ASP:O	3:E:126:ASN:ND2	2.45	0.49
3:E:172:GLY:O	3:E:178:ARG:HA	2.13	0.49
3:H:139:PHE:CB	3:H:193:HIS:HD1	2.26	0.49
1:I:61:ARG:HD3	1:I:66:TYR:CZ	2.47	0.49
1:F:43:ASP:HB3	1:F:45:LYS:H	1.75	0.48
1:F:82:GLU:OE1	2:A:249:TYR:OH	2.24	0.48
2:B:218:ASN:ND2	2:B:221:GLU:HB2	2.27	0.48
1:I:130:LYS:HE2	2:G:247:TYR:HE1	1.77	0.48
2:G:174:ARG:NH1	2:G:178:THR:OG1	2.46	0.48
2:J:376:ALA:HA	2:J:379:THR:HG22	1.95	0.48
1:L:52:PHE:HB2	1:L:55:PHE:H	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:32:ARG:HH12	2:J:36:GLU:HG2	1.77	0.48
2:A:361:PHE:O	3:C:76:LEU:HB2	2.13	0.48
3:E:144:LYS:HE2	3:E:186:ASP:HB2	1.96	0.48
1:I:138:LEU:HD22	2:G:88:TYR:CZ	2.48	0.48
2:A:55:LYS:O	2:A:59:VAL:HG23	2.14	0.48
2:A:66:PHE:HA	2:A:70:HIS:O	2.14	0.48
3:K:90:PHE:CE1	3:K:94:LYS:HE3	2.49	0.48
2:G:232:ARG:O	2:G:236:ILE:HG13	2.14	0.48
2:B:156:VAL:HG22	2:B:191:ARG:CB	2.43	0.48
2:A:96:LEU:O	2:A:100:VAL:HG23	2.14	0.47
3:H:103:SER:HB3	3:H:167:PRO:HB3	1.95	0.47
2:G:192:VAL:O	2:G:195:LEU:HB2	2.14	0.47
3:C:166:HIS:CD2	3:C:167:PRO:HD2	2.49	0.47
1:L:74:ASN:OD1	1:L:74:ASN:N	2.47	0.47
3:H:117:THR:HG22	3:H:118:GLN:HG2	1.95	0.47
2:J:367:LYS:HG2	2:J:396:MET:HB2	1.95	0.47
2:A:107:GLU:OE2	2:A:110:ARG:NH2	2.47	0.47
3:H:64:UNK:C	3:H:66:UNK:HA	2.44	0.47
2:J:147:PHE:O	2:J:151:ILE:HG12	2.14	0.47
2:J:185:MET:HB3	2:J:189:THR:HG21	1.95	0.47
2:A:95:TYR:O	2:A:99:SER:OG	2.13	0.47
1:L:119:LEU:HB3	2:J:170:LEU:HD13	1.97	0.47
1:I:116:GLU:O	1:I:124:ARG:HD3	2.14	0.47
1:L:27:GLU:HG3	1:L:30:LYS:HE3	1.96	0.47
2:G:103:ASN:OD1	2:G:104:SER:N	2.48	0.47
2:G:195:LEU:HA	2:G:195:LEU:HD23	1.80	0.47
1:F:21:MET:HB2	1:F:23:PHE:CE2	2.50	0.47
1:I:42:ARG:NH1	1:I:47:THR:O	2.45	0.47
2:B:100:VAL:HB	2:B:101:LEU:HD12	1.96	0.47
2:J:186:GLY:H	2:J:189:THR:HG23	1.80	0.47
2:B:156:VAL:CG2	2:B:191:ARG:HB2	2.43	0.46
2:B:176:TYR:OH	2:B:217:LYS:HD2	2.15	0.46
2:J:341:ARG:HD2	3:K:177:GLU:OE2	2.15	0.46
2:G:347:TYR:CE2	2:G:388:ARG:HD2	2.50	0.46
2:A:140:SER:OG	2:A:142:GLU:OE1	2.33	0.46
1:L:96:HIS:HB2	3:C:93:GLU:HG2	1.98	0.46
1:I:8:GLN:NE2	1:I:36:HIS:HB2	2.30	0.46
2:G:105:ASN:O	2:G:109:ILE:HG13	2.16	0.46
3:K:74:VAL:HG13	3:K:75:PRO:HD2	1.98	0.46
3:C:136:PRO:HB3	3:C:141:TRP:HD1	1.80	0.46
2:J:293:GLU:HG3	2:J:294:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:K:170:LEU:HD13	3:K:173:MET:HE2	1.98	0.46
2:A:341:ARG:HD2	3:E:177:GLU:OE1	2.16	0.46
3:K:97:LEU:HD23	3:K:97:LEU:HA	1.73	0.46
3:C:97:LEU:HD23	3:C:143:TYR:CE2	2.51	0.46
3:E:117:THR:HG22	3:E:118:GLN:HG2	1.98	0.46
2:A:151:ILE:HB	2:A:152:PRO:HD3	1.97	0.46
2:A:170:LEU:HD23	2:A:170:LEU:HA	1.80	0.45
3:H:89:HIS:O	3:H:93:GLU:HG3	2.16	0.45
3:H:115:TYR:CD1	3:H:122:PRO:HD3	2.51	0.45
1:F:52:PHE:O	1:F:53:ARG:HB3	2.16	0.45
1:I:21:MET:HE2	1:I:23:PHE:CZ	2.51	0.45
1:L:78:LEU:HD23	1:L:78:LEU:HA	1.73	0.45
2:G:123:ARG:HH22	2:G:159:ASP:HB2	1.81	0.45
3:E:115:TYR:CD1	3:E:122:PRO:HD3	2.51	0.45
3:E:121:PHE:CD1	3:E:122:PRO:HD2	2.51	0.45
1:I:34:GLU:O	1:I:38:VAL:HG23	2.16	0.45
2:B:16:PHE:CE1	2:B:36:GLU:HG3	2.51	0.45
2:J:105:ASN:O	2:J:109:ILE:HG13	2.16	0.45
1:L:30:LYS:O	1:L:34:GLU:HG3	2.16	0.45
2:J:193:VAL:HG11	2:J:224:THR:HB	1.98	0.45
3:E:56:UNK:CB	3:E:96:GLY:HA3	2.47	0.45
3:C:165:LEU:HD23	3:C:165:LEU:HA	1.73	0.45
1:I:56:LYS:HE3	1:I:73:VAL:O	2.16	0.45
2:G:170:LEU:HD23	2:G:170:LEU:HA	1.75	0.45
2:A:108:LEU:HD23	2:A:108:LEU:HA	1.78	0.45
3:H:100:LEU:HD23	3:H:183:TRP:CZ3	2.51	0.45
2:G:47:ASP:OD1	2:G:48:LYS:N	2.48	0.45
2:J:117:LYS:HG3	2:J:147:PHE:CE1	2.51	0.45
3:C:166:HIS:CG	3:C:167:PRO:HD2	2.51	0.45
1:F:51:GLU:OE1	2:A:388:ARG:NH2	2.50	0.45
2:B:96:LEU:O	2:B:100:VAL:HG23	2.16	0.45
3:H:147:PRO:HA	3:H:180:VAL:O	2.17	0.45
3:C:111:ASP:HB3	3:C:122:PRO:HB3	1.99	0.45
2:G:370:ILE:HG13	2:G:396:MET:HE3	1.99	0.45
2:B:324:PRO:HB3	2:B:361:PHE:CZ	2.52	0.45
2:J:212:THR:O	2:J:216:GLN:HG3	2.17	0.45
2:A:13:LEU:O	2:A:17:ILE:HG13	2.17	0.44
2:J:288:LEU:HD13	2:J:329:ARG:HB3	1.98	0.44
3:C:97:LEU:HD23	3:C:143:TYR:CZ	2.52	0.44
2:B:129:GLY:O	2:B:133:HIS:ND1	2.50	0.44
2:B:170:LEU:HD23	2:B:170:LEU:HA	1.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:296:LYS:HE2	2:A:285:GLU:OE2	2.18	0.44
2:G:280:LEU:O	2:G:283:CYS:HB3	2.17	0.44
3:E:74:VAL:HG22	3:E:75:PRO:HD2	1.98	0.44
1:F:28:LYS:O	1:F:32:ILE:HG13	2.17	0.44
2:J:204:VAL:O	2:J:208:THR:OG1	2.29	0.44
2:A:324:PRO:HB3	2:A:361:PHE:CZ	2.53	0.44
2:J:377:LEU:HD12	2:J:377:LEU:HA	1.72	0.44
2:A:33:ILE:HG23	2:A:65:ILE:HD13	1.99	0.44
2:A:116:ILE:HG23	2:A:131:ALA:HB1	1.99	0.44
1:F:43:ASP:CG	1:F:44:ALA:HB3	2.38	0.44
2:B:182:LEU:O	2:B:184:PRO:HD3	2.18	0.44
1:D:115:ASP:HB3	2:B:133:HIS:HD2	1.82	0.44
3:K:95:GLY:HA2	3:K:96:GLY:HA2	1.47	0.44
3:K:110:LEU:O	3:K:114:ILE:HG12	2.18	0.44
2:B:13:LEU:O	2:B:17:ILE:HG13	2.17	0.43
2:J:272:GLU:O	2:J:274:PRO:HD3	2.18	0.43
2:B:338:LEU:HD11	2:B:353:MET:HE1	1.99	0.43
3:C:108:ASP:O	3:C:112:LEU:HD23	2.18	0.43
1:F:85:HIS:NE2	1:F:89:GLU:OE2	2.52	0.43
1:L:24:ASP:OD1	1:L:27:GLU:HB3	2.18	0.43
2:J:209:SER:HG	2:J:260:LYS:NZ	2.16	0.43
3:H:100:LEU:HD23	3:H:183:TRP:CE3	2.54	0.43
3:H:196:ARG:HA	3:H:203:PHE:CE2	2.53	0.43
2:G:61:LYS:O	2:G:65:ILE:HG13	2.18	0.43
1:I:31:LEU:HD23	1:I:31:LEU:HA	1.78	0.43
2:J:52:GLY:HA3	2:J:90:GLU:OE2	2.18	0.43
2:A:26:LYS:O	2:A:30:ILE:HD12	2.18	0.43
3:H:166:HIS:CD2	3:H:168:VAL:H	2.37	0.43
2:A:128:MET:HE2	2:A:164:VAL:HG11	1.99	0.43
3:K:152:LYS:HA	3:K:152:LYS:HD3	1.87	0.43
1:I:45:LYS:HB3	1:I:45:LYS:HE2	1.73	0.43
2:J:172:LEU:HD23	2:J:210:LEU:HD21	2.01	0.43
1:L:132:LEU:O	1:L:136:LEU:HD23	2.19	0.43
2:B:212:THR:O	2:B:216:GLN:HG3	2.19	0.43
2:J:91:LYS:NZ	2:J:119:ASP:OD2	2.35	0.43
2:J:96:LEU:O	2:J:100:VAL:HG23	2.19	0.43
3:E:96:GLY:O	3:E:97:LEU:HB2	2.18	0.43
1:L:84:ILE:O	1:L:88:VAL:HG23	2.19	0.43
1:I:43:ASP:OD2	1:I:45:LYS:HE3	2.19	0.43
2:G:241:SER:HB3	2:G:244:LEU:HD13	2.00	0.43
2:G:341:ARG:HD2	3:H:177:GLU:OE2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:11:ARG:O	2:B:15:VAL:HG23	2.19	0.42
2:B:271:PRO:O	2:B:277:ARG:NH2	2.51	0.42
1:F:110:VAL:O	1:F:114:VAL:HG23	2.19	0.42
1:F:12:GLY:HA2	1:F:36:HIS:CD2	2.54	0.42
1:F:127:SER:O	1:F:131:VAL:HG23	2.20	0.42
2:A:189:THR:O	2:A:192:VAL:HG22	2.19	0.42
1:I:87:PHE:CZ	1:I:114:VAL:HG22	2.53	0.42
2:G:83:LEU:O	2:G:91:LYS:HE3	2.19	0.42
3:E:115:TYR:CD1	3:E:121:PHE:HA	2.54	0.42
3:H:125:GLN:HG2	3:H:127:TYR:HE1	1.85	0.42
2:B:282:GLU:OE1	2:A:241:SER:HB3	2.20	0.42
3:H:97:LEU:HD11	3:H:110:LEU:HD21	2.01	0.42
1:I:9:ASN:OD1	1:I:13:LYS:N	2.52	0.42
1:I:60:ARG:NH1	1:I:85:HIS:HB2	2.35	0.42
2:B:317:ILE:HG22	2:B:327:LEU:HD23	2.02	0.42
1:F:98:VAL:CG1	3:E:164:LEU:HD22	2.48	0.42
1:F:138:LEU:HA	1:F:141:LEU:HD22	2.01	0.42
2:B:55:LYS:O	2:B:59:VAL:HG23	2.19	0.42
2:J:322:SER:OG	2:J:323:GLU:HG2	2.20	0.42
2:A:214:LEU:HD23	2:A:214:LEU:HA	1.81	0.42
3:E:196:ARG:HA	3:E:203:PHE:CE2	2.55	0.42
2:J:284:LEU:HD23	2:J:284:LEU:HA	1.78	0.42
2:J:378:LYS:HE2	2:J:378:LYS:HB3	1.87	0.42
3:E:105:ARG:HH22	3:E:167:PRO:HG3	1.85	0.42
1:L:31:LEU:O	1:L:35:VAL:HG23	2.19	0.42
1:I:21:MET:HE3	1:I:21:MET:HB3	1.95	0.42
2:B:83:LEU:O	2:B:91:LYS:HE3	2.20	0.42
2:B:196:LEU:HD12	2:B:196:LEU:HA	1.80	0.42
1:I:43:ASP:OD1	1:I:44:ALA:N	2.50	0.42
1:L:8:GLN:NE2	1:L:32:ILE:HG23	2.34	0.41
2:B:116:ILE:HG23	2:B:131:ALA:HB1	2.02	0.41
2:B:147:PHE:O	2:B:151:ILE:HG12	2.19	0.41
2:G:338:LEU:HD23	2:G:338:LEU:HA	1.87	0.41
2:B:377:LEU:HD12	2:B:377:LEU:HA	1.66	0.41
2:J:151:ILE:HB	2:J:152:PRO:HD3	2.02	0.41
3:H:139:PHE:HB2	3:H:193:HIS:ND1	2.34	0.41
2:G:362:SER:O	2:G:366:VAL:HG23	2.20	0.41
2:J:82:LEU:HD23	2:J:82:LEU:HA	1.89	0.41
2:A:103:ASN:OD1	2:A:104:SER:N	2.53	0.41
1:L:51:GLU:HG3	2:J:347:TYR:OH	2.21	0.41
2:G:301:GLN:HE21	3:H:173:MET:CE	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:123:ARG:NH2	2:J:159:ASP:OD2	2.51	0.41
1:L:30:LYS:HE3	1:L:30:LYS:HB3	1.83	0.41
2:G:125:PRO:O	2:G:128:MET:N	2.53	0.41
2:J:265:LEU:HA	2:J:265:LEU:HD23	1.84	0.41
1:D:15:ARG:HA	1:D:15:ARG:HD2	1.71	0.41
2:B:61:LYS:O	2:B:65:ILE:HG13	2.20	0.41
2:B:158:GLY:O	2:B:159:ASP:OD1	2.39	0.41
3:H:141:TRP:NE1	3:H:143:TYR:HB2	2.35	0.41
3:C:121:PHE:HA	3:C:122:PRO:HD3	1.91	0.41
1:F:90:VAL:CG2	1:F:128:GLN:HG2	2.51	0.41
3:H:104:GLN:N	3:H:174:ASP:OD2	2.23	0.41
3:C:74:VAL:HA	3:C:75:PRO:HD3	1.94	0.41
1:D:38:VAL:CG1	1:D:50:VAL:HG11	2.48	0.41
1:D:52:PHE:HB2	1:D:55:PHE:O	2.20	0.41
1:I:78:LEU:HD23	1:I:78:LEU:HA	1.78	0.41
2:G:67:LEU:HD23	2:G:67:LEU:HA	1.90	0.41
2:B:108:LEU:HD23	2:B:108:LEU:HA	1.92	0.41
2:A:212:THR:O	2:A:216:GLN:HG2	2.21	0.41
1:F:8:GLN:HE21	1:F:8:GLN:HB3	1.68	0.41
1:I:50:VAL:HB	1:I:57:ILE:HB	2.03	0.41
2:G:68:LEU:HD23	2:G:68:LEU:HA	1.90	0.41
2:G:132:LEU:HD23	2:G:132:LEU:HA	1.95	0.41
2:B:31:LYS:O	2:B:35:LYS:HG3	2.20	0.41
2:B:47:ASP:OD1	2:B:48:LYS:N	2.54	0.41
2:B:193:VAL:O	2:B:196:LEU:HB2	2.21	0.41
2:J:294:PRO:HA	2:J:295:PRO:HD3	1.82	0.41
3:H:83:ALA:O	3:H:86:ASP:HB2	2.21	0.41
1:L:76:ASN:N	2:J:263:ARG:HH22	2.19	0.41
1:F:31:LEU:O	1:F:35:VAL:HG23	2.20	0.41
1:I:3:ARG:HG3	1:I:55:PHE:CE2	2.56	0.41
1:I:76:ASN:ND2	2:G:312:GLU:OE1	2.48	0.41
2:B:32:ARG:HH12	2:B:36:GLU:HG2	1.85	0.41
2:J:201:LEU:HA	2:J:201:LEU:HD23	1.75	0.41
3:H:136:PRO:HB3	3:H:141:TRP:HD1	1.84	0.41
1:I:35:VAL:O	1:I:39:VAL:HB	2.21	0.40
2:A:182:LEU:O	2:A:184:PRO:HD3	2.21	0.40
3:H:121:PHE:HA	3:H:122:PRO:HD3	1.88	0.40
1:D:3:ARG:HD2	1:D:21:MET:SD	2.61	0.40
1:L:43:ASP:OD1	1:L:44:ALA:N	2.48	0.40
1:I:3:ARG:HG3	1:I:55:PHE:CZ	2.56	0.40
2:J:124:ASN:O	2:J:128:MET:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:76:ASN:H	2:J:263:ARG:NH2	2.19	0.40
2:J:223:LYS:O	2:J:226:VAL:HG23	2.22	0.40
3:E:77:ARG:N	3:E:119:GLY:O	2.51	0.40
1:F:132:LEU:HD23	1:F:132:LEU:HA	1.97	0.40
2:J:108:LEU:HA	2:J:108:LEU:HD23	1.78	0.40
2:J:123:ARG:HH22	2:J:159:ASP:CG	2.25	0.40
2:J:176:TYR:CE2	2:J:214:LEU:HD23	2.57	0.40
2:A:125:PRO:C	2:A:127:PHE:N	2.74	0.40
3:C:74:VAL:HG13	3:C:75:PRO:HD2	2.03	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	D	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	22	52
1	F	140/142 (99%)	135 (96%)	4 (3%)	1 (1%)	22	52
1	I	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	22	52
1	L	140/142 (99%)	137 (98%)	2 (1%)	1 (1%)	22	52
2	A	381/398 (96%)	378 (99%)	2 (0%)	1 (0%)	41	70
2	B	386/398 (97%)	383 (99%)	3 (1%)	0	100	100
2	G	386/398 (97%)	376 (97%)	6 (2%)	4 (1%)	15	42
2	J	386/398 (97%)	380 (98%)	5 (1%)	1 (0%)	41	70
3	C	130/155 (84%)	128 (98%)	1 (1%)	1 (1%)	19	48
3	E	130/155 (84%)	128 (98%)	2 (2%)	0	100	100
3	H	128/155 (83%)	128 (100%)	0	0	100	100
3	K	129/155 (83%)	128 (99%)	1 (1%)	0	100	100
All	All	2616/2780 (94%)	2575 (98%)	30 (1%)	11 (0%)	34	64

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	159	ASP
1	L	96	HIS
2	G	46	GLY
1	D	96	HIS
1	F	96	HIS
1	I	96	HIS
2	G	158	GLY
3	C	97	LEU
2	G	125	PRO
2	J	46	GLY
2	A	125	PRO

5.3.2 Protein sidechains (i)

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	D	131/131 (100%)	127 (97%)	4 (3%)	40	72
1	F	131/131 (100%)	128 (98%)	3 (2%)	50	79
1	I	131/131 (100%)	123 (94%)	8 (6%)	18	45
1	L	131/131 (100%)	127 (97%)	4 (3%)	40	72
2	A	338/346 (98%)	333 (98%)	5 (2%)	65	86
2	B	340/346 (98%)	334 (98%)	6 (2%)	59	83
2	G	340/346 (98%)	337 (99%)	3 (1%)	78	92
2	J	340/346 (98%)	334 (98%)	6 (2%)	59	83
3	C	118/126 (94%)	113 (96%)	5 (4%)	30	61
3	E	118/126 (94%)	116 (98%)	2 (2%)	60	84
3	H	116/126 (92%)	113 (97%)	3 (3%)	46	76
3	K	117/126 (93%)	114 (97%)	3 (3%)	46	76
All	All	2351/2412 (98%)	2299 (98%)	52 (2%)	52	80

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	15	ARG
1	D	49	PHE
1	D	56	LYS
1	D	136	LEU
1	L	10	ARG
1	L	49	PHE
1	L	116	GLU
1	L	117	MET
1	F	49	PHE
1	F	137	MET
1	F	140	SER
1	I	15	ARG
1	I	16	LEU
1	I	18	LYS
1	I	39	VAL
1	I	49	PHE
1	I	52	PHE
1	I	56	LYS
1	I	82	GLU
2	G	10	MET
2	G	60	CYS
2	G	243	ASP
2	B	10	MET
2	B	60	CYS
2	B	92	GLN
2	B	156	VAL
2	B	159	ASP
2	B	196	LEU
2	J	10	MET
2	J	93	ILE
2	J	104	SER
2	J	293	GLU
2	J	296	LYS
2	J	356	LEU
2	A	91	LYS
2	A	93	ILE
2	A	232	ARG
2	A	246	ASP
2	A	338	LEU
3	E	164	LEU
3	E	183	TRP
3	H	179	GLU
3	H	183	TRP

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Mol	Chain	Res	Type
3	H	188	ARG
3	K	79	MET
3	K	177	GLU
3	K	183	TRP
3	C	76	LEU
3	C	98	GLU
3	C	146	VAL
3	C	183	TRP
3	C	192	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	77	ASN
1	F	8	GLN
2	G	301	GLN
2	B	218	ASN
3	H	125	GLN
3	H	166	HIS
3	K	125	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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	D	142/142 (100%)	-0.18	2 (1%)	75	75	18, 27, 52, 92	0
1	F	142/142 (100%)	-0.06	1 (0%)	87	87	27, 37, 67, 98	0
1	I	142/142 (100%)	-0.21	0	100	100	15, 24, 53, 81	0
1	L	142/142 (100%)	-0.14	1 (0%)	87	87	20, 30, 69, 104	0
2	A	385/398 (96%)	0.10	10 (2%)	56	53	26, 45, 70, 96	0
2	B	388/398 (97%)	-0.02	6 (1%)	73	73	23, 36, 56, 87	0
2	G	388/398 (97%)	0.00	9 (2%)	60	59	20, 34, 55, 78	0
2	J	388/398 (97%)	0.07	8 (2%)	63	62	24, 38, 63, 88	0
3	C	132/155 (85%)	0.19	7 (5%)	26	22	23, 36, 61, 77	0
3	E	132/155 (85%)	0.20	1 (0%)	86	86	19, 37, 56, 69	0
3	H	130/155 (83%)	0.73	21 (16%)	1	1	24, 45, 76, 92	0
3	K	131/155 (84%)	0.41	9 (6%)	16	13	22, 39, 72, 89	0
All	All	2642/2780 (95%)	0.07	75 (2%)	53	50	15, 36, 66, 104	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	47	ASP	5.3
3	H	198	LEU	5.3
2	J	49	ALA	4.4
2	B	47	ASP	4.1
3	H	76	LEU	4.1
3	H	119	GLY	4.0
3	K	75	PRO	4.0
3	H	120	TYR	3.9
3	H	75	PRO	3.8
3	K	73	GLN	3.7
2	A	103	ASN	3.6

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Mol	Chain	Res	Type	RSRZ
3	H	139	PHE	3.5
2	J	48	LYS	3.5
1	F	44	ALA	3.5
2	G	47	ASP	3.5
2	J	46	GLY	3.5
3	C	198	LEU	3.5
3	K	76	LEU	3.3
3	H	80	THR	3.2
3	H	192	HIS	3.2
3	H	203	PHE	3.2
1	L	44	ALA	3.1
3	C	72	PRO	3.1
1	D	44	ALA	3.0
2	J	50	LEU	3.0
2	B	49	ALA	2.9
2	J	25	SER	2.8
3	H	202	TYR	2.8
1	D	53	ARG	2.7
2	B	183	VAL	2.7
3	H	121	PHE	2.7
3	C	194	VAL	2.7
3	H	82	LYS	2.6
3	H	196	ARG	2.6
2	G	394	TYR	2.6
2	A	22	ASN	2.6
3	C	113	TRP	2.6
3	K	116	HIS	2.6
2	B	107	GLU	2.6
2	A	45	LYS	2.5
2	G	42	SER	2.5
3	H	142	CYS	2.5
2	G	46	GLY	2.5
3	K	81	TYR	2.5
3	C	202	TYR	2.5
2	A	29	GLU	2.4
2	B	48	LYS	2.4
3	K	192	HIS	2.3
2	B	185	MET	2.3
3	C	197	GLU	2.3
3	H	118	GLN	2.3
2	A	11	ARG	2.3
3	H	74	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
3	K	80	THR	2.3
3	H	199	HIS	2.3
2	G	123	ARG	2.2
2	G	48	LYS	2.2
3	H	195	ALA	2.2
3	K	74	VAL	2.2
2	A	44	PHE	2.2
2	A	185	MET	2.1
3	K	79	MET	2.1
2	G	360	GLU	2.1
3	C	192	HIS	2.1
3	E	137	LEU	2.1
2	J	44	PHE	2.1
2	G	45	LYS	2.1
3	H	194	VAL	2.1
3	H	193	HIS	2.1
2	G	108	LEU	2.0
2	A	106	SER	2.0
2	A	40	ILE	2.0
2	J	39	ASN	2.0
2	A	104	SER	2.0
3	H	137	LEU	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.