



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 16, 2026 – 02:40 AM JST

PDB ID : 25QB / pdb_000025qb
Title : Topbp1 BRCT0-2 in complex with phosphorylated HTATSF1
Authors : Wang, X.D.; Yang, H.W.; Yang, N.
Deposited on : 2026-04-14
Resolution : 2.99 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:

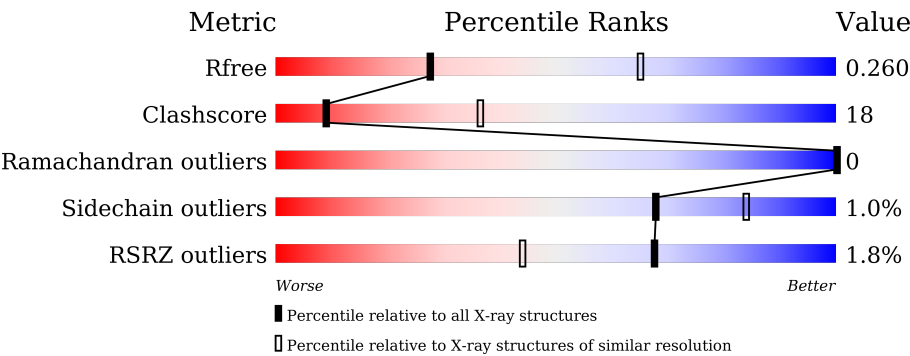
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 2.99 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	281	70%30%
1	B	281	%67%31%..
1	C	281	2%69%28%..
1	D	281	71%28%
1	E	281	4%69%27%. .
1	F	281	2%65%31%..

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Mol	Chain	Length	Quality of chain
2	G	8	88% 12%
2	H	8	88% 12%
2	I	8	75% 25%
2	J	8	12% 75% 25%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13830 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 DNA topoisomerase 2-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	0	0
			2271	1454	382	413	22			
1	B	278	Total	C	N	O	S	0	0	0
			2220	1424	369	405	22			
1	C	276	Total	C	N	O	S	0	0	0
			2219	1423	371	403	22			
1	D	280	Total	C	N	O	S	0	0	0
			2245	1437	376	410	22			
1	E	273	Total	C	N	O	S	0	0	0
			2168	1390	364	392	22			
1	F	275	Total	C	N	O	S	0	0	0
			2208	1418	366	402	22			

- Molecule 2 is a protein called 17S U2 SnRNP complex component HTATSF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	8	Total	C	N	O	P	0	0	0
			58	33	8	16	1			
2	H	8	Total	C	N	O	P	0	0	0
			58	33	8	16	1			
2	I	8	Total	C	N	O	P	0	0	0
			58	33	8	16	1			
2	J	8	Total	C	N	O	P	0	0	0
			58	33	8	16	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	58	Total	O	0	0
			58	58		

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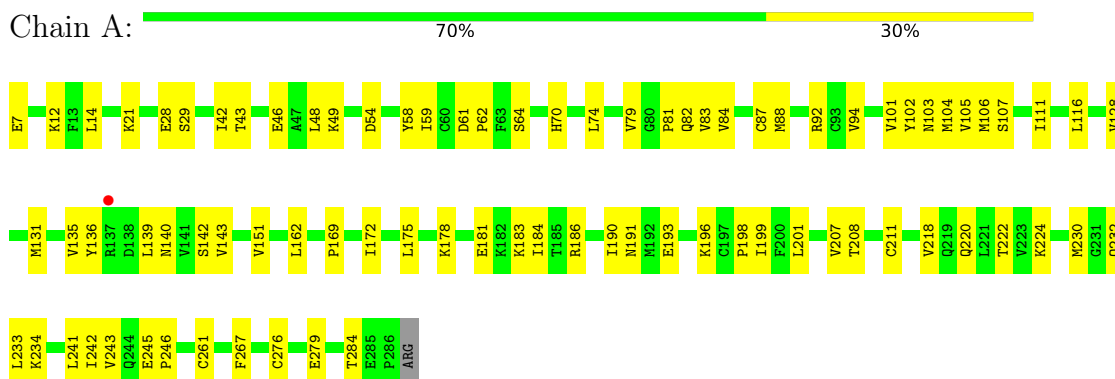
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	40	Total 40	O 40	0	0
3	D	47	Total 47	O 47	0	0
3	E	45	Total 45	O 45	0	0
3	F	28	Total 28	O 28	0	0
3	G	5	Total 5	O 5	0	0
3	J	3	Total 3	O 3	0	0

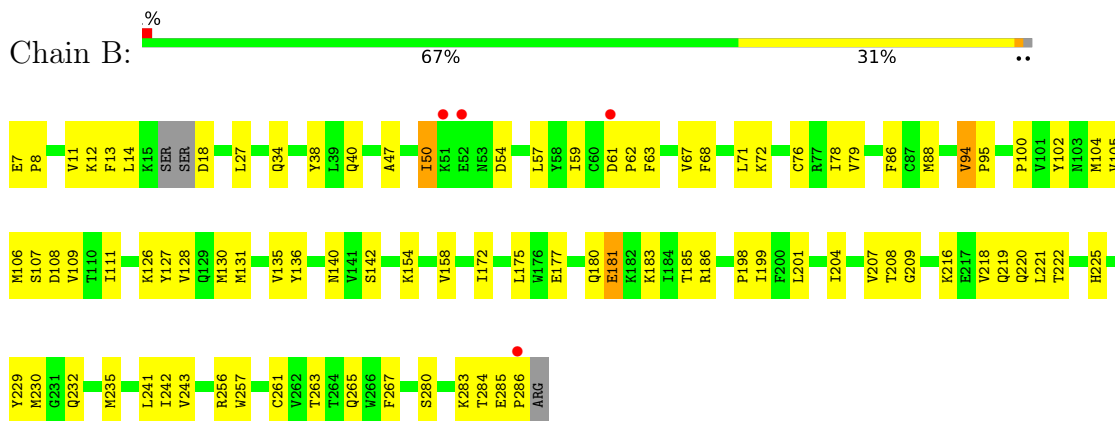
3 Residue-property plots [i](#)

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.

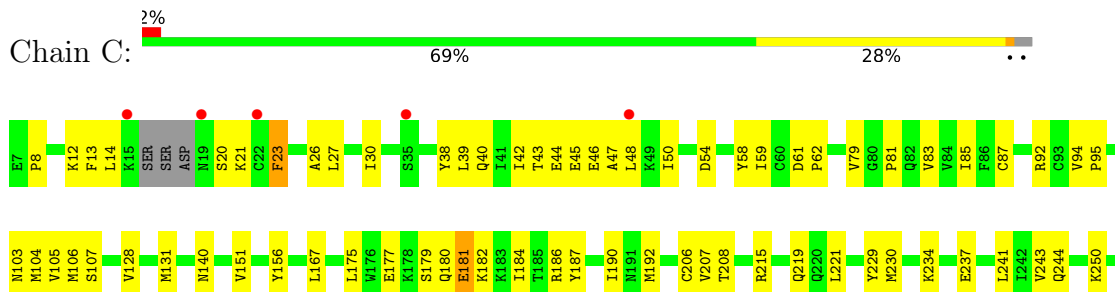
• Molecule 1: DNA topoisomerase 2-binding protein 1

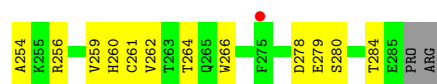


• Molecule 1: DNA topoisomerase 2-binding protein 1



• Molecule 1: DNA topoisomerase 2-binding protein 1





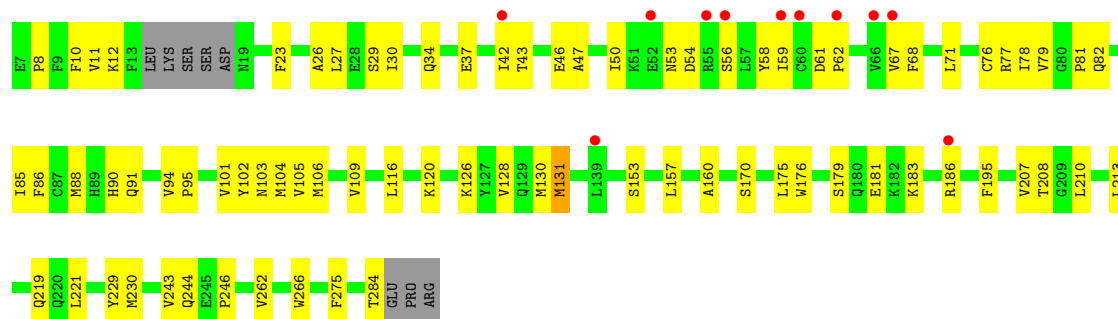
- Molecule 1: DNA topoisomerase 2-binding protein 1

Chain D: 71% 28%



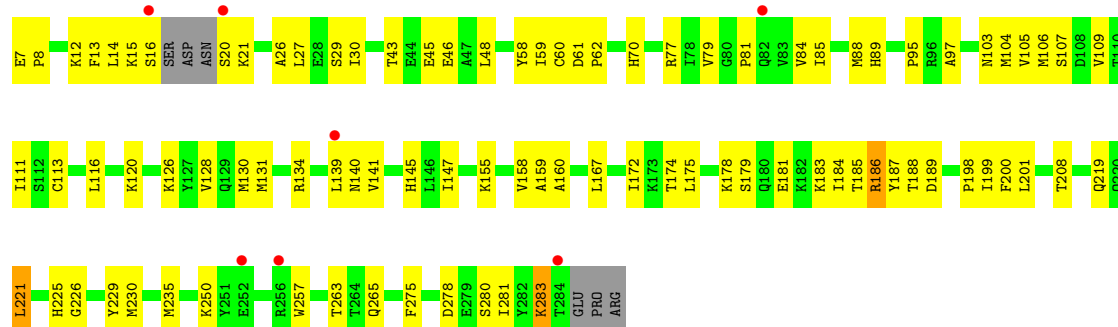
- Molecule 1: DNA topoisomerase 2-binding protein 1

Chain E: 4% 69% 27%



- Molecule 1: DNA topoisomerase 2-binding protein 1

Chain F: 2% 65% 31%



- Molecule 2: 17S U2 SnRNP complex component HTATSF1

Chain G: 88% 12%



- Molecule 2: 17S U2 SnRNP complex component HTATSF1

Chain H:  88% 12%



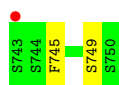
- Molecule 2: 17S U2 SnRNP complex component HTATSF1

Chain I:  75% 25%



- Molecule 2: 17S U2 SnRNP complex component HTATSF1

Chain J:  12% 75% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	123.24Å 163.79Å 170.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.29 – 2.99 40.29 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.29-2.99) 99.8 (40.29-2.99)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.215 , 0.260 0.217 , 0.260	Depositor DCC
R_{free} test set	3539 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	83.0	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13830	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.57	0/2321	0.99	0/3124
1	B	0.65	0/2269	0.98	0/3062
1	C	0.69	0/2267	0.99	0/3055
1	D	0.54	0/2293	0.92	0/3089
1	E	0.64	0/2215	0.96	0/2987
1	F	0.55	0/2255	0.96	0/3037
2	G	0.65	0/47	0.79	0/60
2	H	0.47	0/47	0.57	0/60
2	I	0.60	0/47	0.69	0/60
2	J	0.41	0/47	0.52	0/60
All	All	0.61	0/13808	0.96	0/18594

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

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	2271	0	2278	86	0
1	B	2220	0	2184	90	1
1	C	2219	0	2199	78	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2245	0	2228	94	0
1	E	2168	0	2133	84	0
1	F	2208	0	2200	79	0
2	G	58	0	47	0	0
2	H	58	0	48	0	0
2	I	58	0	48	0	0
2	J	58	0	47	2	0
3	A	41	0	0	6	1
3	B	58	0	0	7	0
3	C	40	0	0	6	0
3	D	47	0	0	5	0
3	E	45	0	0	8	0
3	F	28	0	0	2	0
3	G	5	0	0	0	0
3	J	3	0	0	0	0
All	All	13830	0	13412	497	1

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 18.

All (497) 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:90:HIS:CE1	1:E:186:ARG:NH1	1.68	1.61
1:E:90:HIS:NE2	1:E:186:ARG:NH1	1.61	1.47
1:D:74:LEU:O	1:D:74:LEU:HD23	1.27	1.24
1:B:256:ARG:NH2	1:D:37:GLU:OE1	1.73	1.22
1:E:131:MET:O	3:E:301:HOH:O	1.57	1.18
1:E:54:ASP:OD1	1:E:58:TYR:OH	1.65	1.10
1:D:199:ILE:HG22	1:D:225:HIS:HB3	1.46	0.98
1:A:218:VAL:O	1:A:222:THR:HG23	1.64	0.97
1:B:256:ARG:HH22	1:D:37:GLU:CD	1.71	0.97
1:A:43:THR:HG22	1:A:46:GLU:HG3	1.46	0.96
1:D:53:ASN:HB2	1:D:74:LEU:HD22	1.47	0.95
1:B:47:ALA:O	1:B:50:ILE:HG13	1.68	0.94
1:D:16:SER:C	1:D:18:ASP:N	2.27	0.92
1:F:139:LEU:HD12	1:F:155:LYS:CG	1.99	0.91
1:F:139:LEU:HD12	1:F:155:LYS:HG2	1.52	0.91
1:C:243:VAL:HG23	1:C:284:THR:HG21	1.52	0.91
1:D:53:ASN:HB2	1:D:74:LEU:CD2	2.00	0.90
1:E:103:ASN:N	3:E:303:HOH:O	2.04	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:74:LEU:HD23	1:D:74:LEU:C	1.97	0.88
1:F:184:ILE:CD1	1:F:189:ASP:HB2	2.04	0.87
1:D:74:LEU:O	1:D:74:LEU:CD2	2.18	0.87
1:B:222:THR:HG22	1:B:267:PHE:HE1	1.39	0.86
1:A:233:LEU:O	3:A:301:HOH:O	1.95	0.85
1:E:186:ARG:NH2	3:E:302:HOH:O	2.02	0.83
1:D:219:GLN:HA	1:D:229:TYR:CD2	2.14	0.83
1:A:43:THR:CG2	1:A:46:GLU:HG3	2.11	0.80
1:F:106:MET:HE2	1:F:109:VAL:HG21	1.63	0.80
1:D:219:GLN:HG3	1:D:229:TYR:CE1	2.17	0.79
1:D:217:GLU:O	1:D:221:LEU:HB2	1.82	0.79
1:F:107:SER:HB3	1:F:186:ARG:HH21	1.47	0.79
1:D:208:THR:HG22	1:D:243:VAL:HG12	1.64	0.78
1:C:107:SER:HB3	1:C:186:ARG:HH12	1.48	0.78
1:B:104:MET:HE3	1:B:107:SER:HB2	1.66	0.78
1:E:181:GLU:OE2	1:E:183:LYS:NZ	2.17	0.78
1:C:131:MET:O	3:C:301:HOH:O	2.00	0.77
1:C:256:ARG:NH1	1:E:37:GLU:OE2	2.18	0.76
1:D:151:VAL:HG13	1:D:156:TYR:CZ	2.20	0.75
1:E:153:SER:O	1:E:157:LEU:HD12	1.87	0.74
1:E:43:THR:HG23	1:E:46:GLU:H	1.51	0.74
1:E:54:ASP:CG	1:E:58:TYR:HH	1.92	0.74
1:D:205:ILE:HD12	1:D:222:THR:HG22	1.69	0.74
1:B:47:ALA:O	1:B:50:ILE:CG1	2.36	0.74
1:D:205:ILE:HD12	1:D:222:THR:CG2	2.18	0.74
1:C:105:VAL:N	3:C:301:HOH:O	2.22	0.73
1:A:107:SER:HB3	1:A:186:ARG:NH2	2.04	0.73
1:F:198:PRO:HD2	1:F:201:LEU:HD22	1.69	0.73
1:A:43:THR:HG22	1:A:46:GLU:CG	2.17	0.73
1:A:131:MET:HE1	1:A:175:LEU:HB3	1.71	0.73
1:C:27:LEU:HG	1:C:27:LEU:O	1.88	0.73
1:C:177:GLU:O	1:C:181:GLU:HG3	1.88	0.72
1:D:52:GLU:C	1:D:74:LEU:HD11	2.13	0.72
1:F:184:ILE:HD12	1:F:189:ASP:HB2	1.70	0.72
1:A:29:SER:HB3	1:A:88:MET:HE1	1.70	0.71
1:F:27:LEU:O	1:F:30:ILE:HG22	1.91	0.71
1:D:92:ARG:HH21	1:D:92:ARG:HG2	1.56	0.70
1:E:131:MET:HB3	3:E:301:HOH:O	1.91	0.70
1:F:81:PRO:O	1:F:85:ILE:HD12	1.91	0.70
1:E:82:GLN:OE1	1:E:82:GLN:N	2.14	0.69
1:C:180:GLN:HA	1:C:180:GLN:OE1	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:THR:HG22	1:A:243:VAL:HG12	1.74	0.69
1:D:74:LEU:CD2	1:D:74:LEU:C	2.61	0.69
1:F:29:SER:HB3	1:F:88:MET:HE1	1.75	0.69
1:A:211:CYS:SG	3:A:302:HOH:O	2.49	0.69
1:E:12:LYS:HD3	1:E:42:ILE:HD11	1.74	0.69
1:E:103:ASN:O	3:E:303:HOH:O	2.10	0.69
1:E:105:VAL:HG22	1:E:131:MET:HG2	1.73	0.69
1:F:219:GLN:HG3	1:F:229:TYR:CD1	2.28	0.69
1:D:217:GLU:OE1	3:D:302:HOH:O	2.11	0.68
1:E:86:PHE:HE1	1:E:186:ARG:HE	1.39	0.67
1:E:101:VAL:HG12	3:E:303:HOH:O	1.94	0.67
1:B:180:GLN:OE1	1:B:180:GLN:HA	1.92	0.67
1:D:251:TYR:OH	1:D:255:LYS:HE3	1.95	0.67
1:C:105:VAL:CG2	1:C:131:MET:HE2	2.24	0.67
1:B:208:THR:OG1	1:D:55:ARG:NH1	2.27	0.66
1:F:140:ASN:OD1	1:F:141:VAL:N	2.29	0.66
1:F:139:LEU:HD12	1:F:155:LYS:HG3	1.75	0.66
1:B:222:THR:HG22	1:B:267:PHE:CE1	2.26	0.66
1:A:208:THR:HG22	1:A:243:VAL:CG1	2.26	0.66
1:C:105:VAL:HG21	1:C:131:MET:HE2	1.78	0.66
1:A:82:GLN:HG3	1:A:103:ASN:CB	2.26	0.65
1:E:11:VAL:HG11	1:E:30:ILE:CD1	2.26	0.65
1:C:177:GLU:O	1:C:181:GLU:CG	2.44	0.65
1:F:77:ARG:HD2	1:F:97:ALA:H	1.61	0.65
1:E:90:HIS:CD2	1:E:186:ARG:NH1	2.60	0.65
1:B:142:SER:OG	1:F:201:LEU:CD1	2.46	0.64
1:B:199:ILE:H	1:B:199:ILE:HD12	1.60	0.64
1:D:109:VAL:HG23	1:D:145:HIS:HD1	1.61	0.64
1:F:139:LEU:HB2	1:F:155:LYS:HD3	1.80	0.64
1:A:92:ARG:NH1	3:A:303:HOH:O	2.30	0.63
1:A:181:GLU:O	1:A:183:LYS:N	2.31	0.63
1:E:126:LYS:O	1:E:130:MET:HG3	1.98	0.63
1:F:185:THR:HG22	1:F:189:ASP:OD2	1.99	0.63
1:E:86:PHE:HE1	1:E:186:ARG:NE	1.96	0.63
1:F:139:LEU:HD13	1:F:158:VAL:CG2	2.29	0.63
1:F:107:SER:HB3	1:F:186:ARG:NH2	2.14	0.63
1:B:126:LYS:HE2	1:B:127:TYR:CE2	2.34	0.63
1:C:20:SER:HB3	1:C:23:PHE:HB3	1.81	0.63
1:C:81:PRO:O	1:C:85:ILE:HD12	1.98	0.62
1:D:199:ILE:HG21	1:D:225:HIS:HD2	1.62	0.62
1:D:214:ASP:O	1:D:218:VAL:HG22	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:20:SER:OG	1:F:21:LYS:N	2.31	0.62
1:A:43:THR:HG23	1:A:46:GLU:H	1.64	0.62
1:B:235:MET:HE2	1:B:257:TRP:NE1	2.14	0.62
1:A:103:ASN:OD1	1:A:105:VAL:HG12	1.99	0.62
1:D:207:VAL:HG11	1:D:218:VAL:HG21	1.81	0.62
1:B:218:VAL:O	1:B:222:THR:HG23	2.00	0.61
1:C:48:LEU:O	1:C:48:LEU:HD23	2.00	0.61
1:E:94:VAL:HG13	1:E:95:PRO:HD2	1.83	0.61
1:A:279:GLU:OE2	1:A:279:GLU:N	2.34	0.61
1:A:82:GLN:HG3	1:A:103:ASN:HB3	1.82	0.60
1:E:246:PRO:HG3	1:E:284:THR:HG22	1.81	0.60
1:B:285:GLU:HG3	1:B:286:PRO:HD2	1.82	0.60
1:D:53:ASN:HB2	1:D:74:LEU:HD21	1.83	0.60
1:B:263:THR:OG1	1:B:265:GLN:HG3	2.02	0.60
1:F:179:SER:HA	1:F:184:ILE:HG22	1.83	0.60
1:B:88:MET:HG3	3:B:350:HOH:O	2.02	0.60
1:C:206:CYS:HB2	1:C:241:LEU:HD23	1.84	0.59
1:C:244:GLN:OE1	1:C:264:THR:HG23	2.02	0.59
1:D:219:GLN:HG3	1:D:229:TYR:CD1	2.37	0.59
1:D:222:THR:CG2	1:D:267:PHE:HE1	2.16	0.59
1:C:43:THR:HG23	1:C:46:GLU:OE1	2.02	0.59
1:F:184:ILE:HD11	1:F:189:ASP:HB2	1.83	0.59
1:F:103:ASN:OD1	1:F:105:VAL:HG13	2.03	0.59
1:B:104:MET:O	1:B:107:SER:HB3	2.03	0.59
1:D:181:GLU:O	1:D:183:LYS:N	2.34	0.59
1:E:170:SER:HG	1:E:195:PHE:HD2	1.50	0.58
1:C:208:THR:HG22	1:C:243:VAL:CG1	2.33	0.58
1:D:208:THR:HG22	1:D:243:VAL:CG1	2.34	0.58
1:C:206:CYS:CB	1:C:241:LEU:HD23	2.34	0.58
1:E:105:VAL:HG21	1:E:131:MET:HE2	1.86	0.58
1:E:11:VAL:HG11	1:E:30:ILE:HD11	1.85	0.58
1:B:40:GLN:NE2	3:B:302:HOH:O	2.37	0.58
1:D:44:GLU:O	1:D:48:LEU:HD22	2.04	0.57
1:F:105:VAL:HG22	1:F:131:MET:HB3	1.85	0.57
1:E:219:GLN:HG3	1:E:229:TYR:CG	2.40	0.57
1:C:42:ILE:HD12	1:C:46:GLU:HB2	1.87	0.57
1:E:47:ALA:HA	1:E:50:ILE:HD12	1.86	0.57
1:F:48:LEU:HD13	1:F:48:LEU:O	2.05	0.57
1:B:105:VAL:O	1:B:106:MET:HB3	2.04	0.57
1:B:86:PHE:CD1	1:B:104:MET:HG2	2.40	0.57
1:D:207:VAL:HG12	1:D:242:ILE:HB	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:GLU:HA	1:A:196:LYS:HG3	1.87	0.56
1:E:56:SER:O	1:E:58:TYR:CD1	2.57	0.56
1:A:241:LEU:HG	1:A:243:VAL:HG13	1.85	0.56
1:D:281:ILE:HG23	1:D:282:TYR:CD2	2.40	0.56
1:C:23:PHE:CD1	1:C:23:PHE:C	2.83	0.56
1:F:199:ILE:HD12	1:F:199:ILE:H	1.70	0.56
1:B:204:ILE:CG2	1:B:230:MET:HG2	2.34	0.56
1:C:184:ILE:HG21	1:C:190:ILE:CD1	2.36	0.56
1:D:281:ILE:HG23	1:D:282:TYR:CE2	2.40	0.56
1:F:77:ARG:NH1	1:F:95:PRO:O	2.38	0.56
1:B:38:TYR:OH	3:B:301:HOH:O	2.10	0.56
1:C:79:VAL:HB	1:C:83:VAL:HG21	1.88	0.56
1:B:68:PHE:HE1	1:B:72:LYS:HE2	1.71	0.55
1:A:7:GLU:N	3:A:305:HOH:O	2.39	0.55
1:D:52:GLU:CA	1:D:74:LEU:HD11	2.37	0.55
1:E:67:VAL:O	1:E:71:LEU:HD12	2.06	0.55
1:F:278:ASP:HB3	1:F:281:ILE:HD13	1.88	0.55
1:E:207:VAL:HB	1:E:210:LEU:HD13	1.88	0.55
1:D:247:LYS:HE3	1:D:248:GLY:H	1.72	0.54
1:B:241:LEU:HG	1:B:243:VAL:HG23	1.89	0.54
1:D:55:ARG:NH2	3:D:301:HOH:O	2.01	0.54
1:D:219:GLN:CG	1:D:229:TYR:CD1	2.90	0.54
1:E:131:MET:HE1	1:E:175:LEU:C	2.33	0.54
1:E:131:MET:HE1	1:E:176:TRP:N	2.23	0.54
1:A:83:VAL:HG21	1:A:101:VAL:HG23	1.90	0.54
1:E:54:ASP:CG	1:E:58:TYR:OH	2.46	0.54
1:E:106:MET:N	3:E:301:HOH:O	1.97	0.54
1:D:174:THR:O	1:D:178:LYS:HB2	2.07	0.54
1:F:7:GLU:N	1:F:8:PRO:HD2	2.22	0.53
1:C:43:THR:OG1	1:C:45:GLU:HB3	2.08	0.53
1:D:222:THR:HG23	1:D:267:PHE:CE1	2.43	0.53
1:F:77:ARG:HG3	1:F:97:ALA:HB3	1.91	0.53
1:C:131:MET:HE1	1:C:175:LEU:C	2.34	0.53
1:B:68:PHE:CE1	1:B:72:LYS:HE2	2.44	0.53
1:B:177:GLU:O	1:B:181:GLU:HG3	2.08	0.53
1:F:26:ALA:HB2	1:F:85:ILE:HD11	1.91	0.53
1:D:229:TYR:O	1:D:230:MET:HE2	2.09	0.53
1:E:219:GLN:HG3	1:E:229:TYR:CD1	2.44	0.53
1:A:29:SER:HB3	1:A:88:MET:CE	2.39	0.53
1:B:11:VAL:HG12	1:B:57:LEU:HB3	1.90	0.53
1:C:26:ALA:HB2	1:C:85:ILE:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:ALA:O	1:D:163:LYS:HE2	2.08	0.53
1:E:106:MET:HG2	1:E:131:MET:HB3	1.90	0.53
1:A:234:LYS:HA	3:A:301:HOH:O	2.08	0.52
1:C:254:ALA:HB1	1:C:259:VAL:HG22	1.90	0.52
1:A:198:PRO:HB2	1:A:201:LEU:HB2	1.91	0.52
1:A:207:VAL:HG11	1:A:218:VAL:HG21	1.91	0.52
1:D:31:LYS:NZ	3:D:306:HOH:O	2.42	0.52
1:A:178:LYS:HD2	1:A:190:ILE:HD11	1.90	0.52
1:A:222:THR:HG22	1:A:267:PHE:CE1	2.44	0.52
1:A:199:ILE:HD12	1:A:199:ILE:H	1.75	0.52
1:D:28:GLU:OE2	3:D:303:HOH:O	2.19	0.52
1:F:235:MET:HG3	1:F:257:TRP:CD2	2.45	0.52
1:F:278:ASP:HB3	1:F:281:ILE:CD1	2.40	0.52
1:A:70:HIS:CE1	1:A:74:LEU:HD11	2.45	0.51
1:E:47:ALA:HA	1:E:50:ILE:CD1	2.39	0.51
1:E:81:PRO:O	1:E:85:ILE:HD12	2.09	0.51
1:E:106:MET:SD	1:E:128:VAL:HG23	2.50	0.51
1:B:14:LEU:HD11	1:B:67:VAL:CG2	2.41	0.51
1:C:208:THR:HG22	1:C:243:VAL:HG12	1.93	0.51
1:E:27:LEU:O	1:E:30:ILE:HG22	2.11	0.51
1:C:131:MET:HB3	3:C:301:HOH:O	2.11	0.51
1:A:261:CYS:C	1:A:284:THR:HG23	2.36	0.51
1:F:104:MET:HB3	1:F:107:SER:HB2	1.93	0.51
1:B:241:LEU:HG	1:B:243:VAL:CG2	2.41	0.51
1:F:77:ARG:CG	1:F:97:ALA:HB3	2.41	0.51
1:B:207:VAL:HG12	1:B:242:ILE:HB	1.92	0.50
1:D:176:TRP:O	1:D:179:SER:HB3	2.10	0.50
1:B:111:ILE:HG22	1:B:135:VAL:HA	1.93	0.50
1:C:14:LEU:HD12	1:C:42:ILE:HG23	1.94	0.50
1:A:21:LYS:HD2	1:A:21:LYS:O	2.11	0.50
1:B:13:PHE:CE1	1:B:59:ILE:HD12	2.46	0.50
1:D:53:ASN:ND2	1:D:55:ARG:H	2.09	0.50
1:A:107:SER:HB3	1:A:186:ARG:HH21	1.76	0.50
1:A:178:LYS:O	1:A:184:ILE:HG12	2.12	0.50
1:B:154:LYS:O	1:B:158:VAL:HG23	2.11	0.50
1:B:199:ILE:HD12	1:B:199:ILE:N	2.27	0.50
1:C:175:LEU:O	1:C:179:SER:HB3	2.11	0.50
1:D:84:VAL:O	1:D:88:MET:HG2	2.12	0.50
1:D:222:THR:HG23	1:D:267:PHE:HE1	1.77	0.50
1:C:13:PHE:O	1:C:42:ILE:HG22	2.12	0.50
1:A:79:VAL:HG13	1:A:83:VAL:HB	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:SER:HB3	1:C:23:PHE:CB	2.41	0.50
1:F:81:PRO:O	1:F:84:VAL:HG22	2.12	0.50
1:B:67:VAL:O	1:B:71:LEU:HD12	2.12	0.49
1:E:12:LYS:HD3	1:E:42:ILE:CD1	2.40	0.49
1:E:47:ALA:O	1:E:50:ILE:HD12	2.12	0.49
1:C:186:ARG:HH21	1:C:187:TYR:HE1	1.60	0.49
1:E:244:GLN:HA	1:E:244:GLN:OE1	2.12	0.49
1:E:77:ARG:NH2	1:E:94:VAL:HG12	2.27	0.49
1:B:263:THR:OG1	1:B:265:GLN:CG	2.59	0.49
1:E:90:HIS:NE2	1:E:186:ARG:CZ	2.63	0.49
1:E:131:MET:HE3	1:E:176:TRP:HA	1.94	0.49
1:A:49:LYS:HD2	1:A:49:LYS:N	2.27	0.49
1:A:230:MET:C	1:A:232:GLN:H	2.20	0.49
1:B:67:VAL:HG12	3:B:337:HOH:O	2.12	0.49
1:D:199:ILE:HG21	1:D:225:HIS:CD2	2.46	0.49
1:E:105:VAL:N	3:E:301:HOH:O	2.45	0.49
1:E:54:ASP:OD1	1:E:58:TYR:CZ	2.61	0.49
1:A:81:PRO:O	1:A:84:VAL:HG22	2.12	0.49
1:F:175:LEU:O	1:F:179:SER:HB3	2.12	0.49
1:B:216:LYS:O	1:B:220:GLN:HG3	2.13	0.49
1:C:54:ASP:HB3	1:C:58:TYR:OH	2.13	0.49
1:D:151:VAL:HG13	1:D:156:TYR:CE1	2.47	0.49
1:B:199:ILE:HD13	1:B:225:HIS:ND1	2.27	0.48
1:C:14:LEU:HA	1:C:42:ILE:HG23	1.95	0.48
1:C:219:GLN:HG3	1:C:229:TYR:CG	2.48	0.48
1:D:216:LYS:O	1:D:220:GLN:HG3	2.13	0.48
1:E:131:MET:CE	1:E:176:TRP:HA	2.43	0.48
1:F:139:LEU:C	1:F:139:LEU:HD23	2.39	0.48
1:F:59:ILE:HA	1:F:79:VAL:O	2.13	0.48
1:F:134:ARG:HD2	3:F:308:HOH:O	2.12	0.48
1:C:254:ALA:HB1	1:C:259:VAL:CG2	2.43	0.48
1:E:77:ARG:HH21	1:E:94:VAL:HG12	1.78	0.48
1:F:208:THR:OG1	1:F:250:LYS:HD2	2.13	0.48
1:B:18:ASP:N	3:B:303:HOH:O	2.46	0.48
1:D:11:VAL:HG12	1:D:57:LEU:HB3	1.95	0.48
1:E:106:MET:HE2	1:E:109:VAL:HG21	1.95	0.48
1:D:205:ILE:HD12	1:D:222:THR:HG21	1.92	0.48
1:A:111:ILE:HD13	1:A:128:VAL:HG11	1.94	0.48
1:B:204:ILE:HG22	1:B:230:MET:HG2	1.96	0.48
1:B:131:MET:HE1	1:B:175:LEU:HB3	1.95	0.48
1:B:136:TYR:HE2	1:B:142:SER:HB2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:281:ILE:H	1:F:281:ILE:HD12	1.79	0.48
1:D:255:LYS:HE2	1:D:255:LYS:HB2	1.47	0.47
1:E:170:SER:OG	1:E:195:PHE:HD2	1.97	0.47
1:D:219:GLN:HG2	1:D:229:TYR:CG	2.48	0.47
1:A:151:VAL:HG11	1:A:276:CYS:CB	2.44	0.47
1:B:105:VAL:O	1:B:131:MET:HB3	2.15	0.47
1:A:79:VAL:CG1	1:A:83:VAL:HB	2.45	0.47
1:B:13:PHE:CD2	1:B:27:LEU:HD13	2.50	0.47
1:B:61:ASP:HB3	1:B:62:PRO:HD3	1.97	0.47
1:E:208:THR:HG22	1:E:243:VAL:HG22	1.96	0.47
1:B:94:VAL:HG13	1:B:95:PRO:HD2	1.97	0.47
1:C:180:GLN:C	1:C:182:LYS:H	2.22	0.47
1:B:221:LEU:HD23	1:B:221:LEU:HA	1.78	0.47
1:F:48:LEU:HD22	1:F:70:HIS:CD2	2.50	0.47
1:F:181:GLU:O	1:F:183:LYS:N	2.47	0.47
1:D:44:GLU:O	1:D:48:LEU:CD2	2.63	0.47
1:B:14:LEU:HD11	1:B:67:VAL:HG22	1.96	0.47
1:B:106:MET:HE1	1:B:172:ILE:CD1	2.45	0.47
1:D:109:VAL:HG23	1:D:144:THR:OG1	2.14	0.47
1:C:61:ASP:HB3	1:C:62:PRO:HD3	1.96	0.47
1:E:76:CYS:O	1:E:78:ILE:HD12	2.15	0.47
1:F:174:THR:O	1:F:178:LYS:HB2	2.15	0.47
1:F:139:LEU:CD1	1:F:155:LYS:HG3	2.41	0.46
1:C:262:VAL:HG21	1:C:266:TRP:CE3	2.50	0.46
1:D:61:ASP:HB3	1:D:62:PRO:HD3	1.97	0.46
1:C:184:ILE:HG21	1:C:190:ILE:HD11	1.97	0.46
1:E:120:LYS:O	1:E:120:LYS:HG3	2.15	0.46
1:C:207:VAL:HG21	1:C:215:ARG:HG2	1.98	0.46
1:C:250:LYS:HG3	1:E:10:PHE:CE1	2.51	0.46
1:D:126:LYS:HE3	1:D:130:MET:HE2	1.97	0.46
1:B:59:ILE:HA	1:B:79:VAL:O	2.15	0.46
1:F:263:THR:HG23	1:F:265:GLN:HB2	1.98	0.46
1:A:12:LYS:HD2	1:A:58:TYR:CE2	2.51	0.46
1:F:281:ILE:HD12	1:F:281:ILE:N	2.30	0.46
1:C:260:HIS:CG	1:C:279:GLU:HG2	2.51	0.46
1:D:53:ASN:N	1:D:74:LEU:HD11	2.31	0.46
1:A:106:MET:CE	1:A:128:VAL:HA	2.46	0.46
1:B:198:PRO:HD2	1:B:201:LEU:HD22	1.98	0.46
1:C:278:ASP:OD2	1:C:280:SER:OG	2.34	0.46
1:B:198:PRO:HB2	1:B:201:LEU:HB2	1.98	0.46
1:B:105:VAL:O	1:B:131:MET:O	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:GLY:H	1:D:55:ARG:HH11	1.64	0.45
1:C:221:LEU:HA	1:C:221:LEU:HD23	1.65	0.45
1:E:82:GLN:NE2	1:E:102:TYR:CE2	2.85	0.45
1:A:43:THR:HG22	1:A:46:GLU:OE2	2.16	0.45
1:B:95:PRO:HG3	1:B:104:MET:HE1	1.97	0.45
1:E:61:ASP:HB3	1:E:62:PRO:HD3	1.99	0.45
1:B:219:GLN:HG2	1:B:229:TYR:CG	2.51	0.45
1:B:256:ARG:NH2	1:D:37:GLU:HG3	2.31	0.45
1:A:87:CYS:SG	1:A:94:VAL:HA	2.57	0.45
1:C:30:ILE:HG21	1:C:39:LEU:HD12	1.99	0.45
1:A:135:VAL:HG22	3:A:318:HOH:O	2.17	0.45
1:C:42:ILE:CD1	1:C:46:GLU:HB2	2.46	0.45
1:D:92:ARG:HG2	1:D:92:ARG:NH2	2.28	0.45
1:A:48:LEU:HA	1:A:70:HIS:CD2	2.52	0.45
1:B:105:VAL:HG22	1:B:185:THR:C	2.42	0.45
1:B:219:GLN:HG2	1:B:229:TYR:CD1	2.51	0.45
1:C:42:ILE:HD12	1:C:46:GLU:CB	2.47	0.45
1:C:87:CYS:HB3	1:C:92:ARG:O	2.17	0.45
1:D:198:PRO:HB2	1:D:201:LEU:HB2	1.97	0.45
1:E:86:PHE:CE1	1:E:186:ARG:NH1	2.84	0.45
1:F:160:ALA:HB1	1:F:275:PHE:CG	2.52	0.45
1:A:106:MET:HB2	1:A:131:MET:HB3	1.99	0.45
1:B:126:LYS:HE2	1:B:127:TYR:CZ	2.50	0.45
1:C:43:THR:HG1	1:C:45:GLU:HB3	1.82	0.45
1:D:178:LYS:HB3	1:D:184:ILE:CD1	2.47	0.45
1:F:219:GLN:HG3	1:F:229:TYR:CG	2.52	0.45
1:A:116:LEU:HD21	1:A:169:PRO:HG3	1.99	0.45
1:A:136:TYR:HE2	1:A:142:SER:HB2	1.82	0.45
1:B:232:GLN:HA	1:B:232:GLN:OE1	2.16	0.45
1:D:120:LYS:HE2	1:D:169:PRO:CB	2.46	0.45
1:F:89:HIS:HD1	1:F:89:HIS:C	2.24	0.45
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.78	0.44
1:D:109:VAL:HG23	1:D:145:HIS:ND1	2.32	0.44
1:E:116:LEU:HD23	1:E:116:LEU:HA	1.82	0.44
1:A:151:VAL:HG11	1:A:276:CYS:HB2	1.99	0.44
1:B:106:MET:HE2	1:B:109:VAL:HG21	2.00	0.44
1:C:140:ASN:OD1	1:C:140:ASN:C	2.61	0.44
1:D:219:GLN:HG3	1:D:229:TYR:CZ	2.52	0.44
1:F:283:LYS:HD2	1:F:283:LYS:HA	1.76	0.44
1:A:105:VAL:HG23	1:A:186:ARG:CA	2.48	0.44
1:C:261:CYS:C	1:C:284:THR:HG23	2.42	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:186:ARG:O	1:F:188:THR:N	2.51	0.44
1:A:61:ASP:HB3	1:A:62:PRO:HD3	1.98	0.44
1:B:76:CYS:O	1:B:78:ILE:HD12	2.16	0.44
1:B:256:ARG:HH22	1:D:37:GLU:CG	2.31	0.44
1:E:91:GLN:O	1:E:91:GLN:HG2	2.18	0.44
1:B:126:LYS:O	1:B:130:MET:HG3	2.18	0.44
1:A:105:VAL:HG11	1:A:131:MET:HE3	1.99	0.44
1:A:107:SER:H	1:A:186:ARG:HH21	1.65	0.44
1:B:257:TRP:CZ2	1:D:37:GLU:HG2	2.53	0.44
1:D:219:GLN:HG2	1:D:229:TYR:CD1	2.53	0.44
1:A:28:GLU:HG3	1:A:29:SER:H	1.83	0.44
1:A:140:ASN:C	1:A:140:ASN:OD1	2.61	0.44
1:A:222:THR:HG22	1:A:267:PHE:HE1	1.83	0.44
1:C:44:GLU:O	3:C:302:HOH:O	2.21	0.44
1:D:55:ARG:NH1	3:D:309:HOH:O	2.51	0.44
1:C:48:LEU:N	3:C:302:HOH:O	2.39	0.43
1:C:167:LEU:HD11	1:C:192:MET:HG2	2.00	0.43
1:D:106:MET:HE2	1:D:106:MET:HB3	1.78	0.43
1:E:59:ILE:HA	1:E:79:VAL:O	2.18	0.43
1:F:113:CYS:HB3	1:F:116:LEU:HD12	2.00	0.43
1:A:21:LYS:HD2	1:A:21:LYS:C	2.43	0.43
1:A:61:ASP:O	1:A:102:TYR:HE1	2.01	0.43
1:B:106:MET:CE	1:B:109:VAL:HG21	2.48	0.43
1:E:230:MET:HE3	1:E:230:MET:HB3	1.73	0.43
1:E:262:VAL:HG21	1:E:266:TRP:CE3	2.53	0.43
1:F:13:PHE:CE1	1:F:59:ILE:HD12	2.53	0.43
1:F:15:LYS:HG3	1:F:16:SER:N	2.33	0.43
1:B:140:ASN:ND2	1:F:226:GLY:HA2	2.34	0.43
1:A:14:LEU:HD22	1:A:42:ILE:O	2.18	0.43
1:D:199:ILE:HG13	1:D:275:PHE:HA	2.00	0.43
1:B:72:LYS:HD3	1:B:100:PRO:HG2	2.00	0.43
1:B:135:VAL:HG22	3:B:330:HOH:O	2.18	0.43
1:E:160:ALA:HB1	1:E:275:PHE:CD1	2.53	0.43
1:B:105:VAL:CG1	1:B:131:MET:HE2	2.49	0.43
1:D:14:LEU:HD12	1:D:15:LYS:N	2.33	0.43
1:D:173:LYS:HE3	1:D:173:LYS:HB2	1.87	0.43
1:A:139:LEU:HD12	1:A:143:VAL:HG21	2.00	0.43
1:B:106:MET:HB3	1:B:131:MET:HB3	2.00	0.43
1:B:106:MET:HG3	1:B:109:VAL:CG2	2.48	0.43
1:C:30:ILE:CG2	1:C:39:LEU:HD12	2.48	0.43
1:D:14:LEU:HD12	1:D:15:LYS:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:VAL:HG12	1:A:242:ILE:HB	2.00	0.43
1:C:8:PRO:HB3	1:C:38:TYR:CE2	2.54	0.43
1:D:90:HIS:O	1:D:91:GLN:HB2	2.19	0.43
1:E:56:SER:O	1:E:58:TYR:CE1	2.72	0.43
1:A:232:GLN:OE1	1:A:232:GLN:HA	2.19	0.42
1:B:256:ARG:CZ	1:D:37:GLU:OE1	2.57	0.42
1:B:12:LYS:NZ	1:B:54:ASP:OD2	2.52	0.42
1:C:94:VAL:HG13	1:C:95:PRO:HD2	2.01	0.42
1:C:234:LYS:HE3	1:C:237:GLU:OE1	2.18	0.42
1:E:88:MET:HE3	1:E:88:MET:HB3	1.86	0.42
1:E:175:LEU:O	1:E:179:SER:HB3	2.19	0.42
1:F:88:MET:HB2	1:F:88:MET:HE2	1.74	0.42
1:F:130:MET:HE2	1:F:130:MET:HB3	1.95	0.42
1:A:207:VAL:CG1	1:A:218:VAL:HG21	2.49	0.42
1:C:12:LYS:HA	1:C:40:GLN:O	2.19	0.42
1:C:219:GLN:HG3	1:C:229:TYR:CD1	2.54	0.42
1:F:139:LEU:HD11	1:F:159:ALA:HB2	2.00	0.42
1:B:7:GLU:N	3:B:305:HOH:O	2.52	0.42
1:D:16:SER:O	1:D:18:ASP:N	2.53	0.42
1:E:104:MET:HE3	1:E:104:MET:HB3	1.85	0.42
1:A:245:GLU:HA	1:A:246:PRO:HD3	1.86	0.42
1:D:199:ILE:HD11	1:D:270:SER:O	2.20	0.42
1:E:23:PHE:HB2	1:E:61:ASP:OD2	2.19	0.42
1:E:85:ILE:HD12	1:E:85:ILE:H	1.85	0.42
1:F:126:LYS:O	1:F:130:MET:HG3	2.19	0.42
1:B:63:PHE:CD1	1:B:102:TYR:HD1	2.37	0.42
1:D:221:LEU:HD23	1:D:221:LEU:HA	1.83	0.42
1:A:106:MET:HE3	1:A:131:MET:HB2	2.01	0.42
1:D:94:VAL:HG13	1:D:95:PRO:HD2	2.01	0.42
1:A:62:PRO:C	1:A:64:SER:H	2.28	0.42
1:E:8:PRO:HA	1:E:34:GLN:OE1	2.20	0.42
1:E:105:VAL:CG2	1:E:131:MET:HE2	2.49	0.42
1:E:213:LEU:HD22	1:E:213:LEU:H	1.85	0.42
1:A:128:VAL:HG12	1:A:172:ILE:HD12	2.02	0.41
1:B:131:MET:HE1	1:B:175:LEU:C	2.45	0.41
1:C:177:GLU:O	1:C:181:GLU:HG2	2.19	0.41
1:D:52:GLU:N	1:D:52:GLU:OE1	2.53	0.41
1:F:27:LEU:HD12	1:F:27:LEU:HA	1.79	0.41
1:F:278:ASP:OD2	1:F:280:SER:OG	2.37	0.41
1:A:79:VAL:HG22	1:A:101:VAL:HG22	2.02	0.41
1:F:14:LEU:HD22	1:F:60:CYS:SG	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:CYS:O	1:A:284:THR:HG23	2.19	0.41
1:C:104:MET:O	1:C:107:SER:HB2	2.20	0.41
1:C:207:VAL:CG2	1:C:215:ARG:HG2	2.50	0.41
1:D:28:GLU:O	1:D:29:SER:C	2.62	0.41
1:E:26:ALA:HB2	1:E:85:ILE:HD11	2.02	0.41
1:F:147:ILE:HA	1:F:167:LEU:O	2.21	0.41
1:A:230:MET:O	1:A:232:GLN:N	2.54	0.41
1:B:128:VAL:HG12	1:B:172:ILE:HD12	2.02	0.41
1:C:106:MET:SD	1:C:128:VAL:HG23	2.60	0.41
1:D:61:ASP:O	1:D:102:TYR:HE1	2.04	0.41
1:D:247:LYS:HA	1:D:247:LYS:HD2	1.79	0.41
1:E:29:SER:C	1:E:88:MET:HE1	2.44	0.41
1:B:8:PRO:HA	1:B:34:GLN:OE1	2.19	0.41
1:B:280:SER:O	1:B:283:LYS:HE2	2.19	0.41
1:C:241:LEU:HD13	1:C:243:VAL:HG13	2.03	0.41
1:D:12:LYS:HA	1:D:40:GLN:O	2.20	0.41
1:F:200:PHE:HD2	1:F:225:HIS:HB2	1.86	0.41
1:F:221:LEU:HD12	1:F:221:LEU:HA	1.82	0.41
1:E:160:ALA:HB1	1:E:275:PHE:CG	2.55	0.41
1:A:175:LEU:O	1:A:184:ILE:HD11	2.21	0.41
1:C:107:SER:CB	1:C:186:ARG:HH12	2.23	0.41
1:D:118:LYS:H	1:D:118:LYS:HG3	1.61	0.41
1:F:43:THR:OG1	1:F:45:GLU:HG3	2.21	0.41
1:F:60:CYS:O	1:F:81:PRO:HD3	2.20	0.41
1:A:104:MET:HB3	1:A:107:SER:HB2	2.03	0.41
1:A:105:VAL:HG23	1:A:186:ARG:C	2.45	0.41
1:A:131:MET:HE1	1:A:175:LEU:CB	2.48	0.41
1:B:108:ASP:H	1:B:186:ARG:HH22	1.67	0.41
1:C:104:MET:C	3:C:301:HOH:O	2.62	0.41
1:C:151:VAL:HG23	1:C:156:TYR:CE2	2.56	0.41
1:C:230:MET:SD	1:C:237:GLU:CD	3.03	0.41
1:D:208:THR:OG1	1:D:250:LYS:HD2	2.20	0.41
1:F:111:ILE:HG22	1:F:145:HIS:HB2	2.03	0.41
1:A:224:LYS:HB2	2:J:745:PHE:CE1	2.56	0.41
1:B:12:LYS:HA	1:B:40:GLN:O	2.21	0.41
1:B:95:PRO:HG3	1:B:104:MET:SD	2.61	0.41
1:C:47:ALA:HA	1:C:50:ILE:CD1	2.51	0.41
1:C:243:VAL:O	1:C:284:THR:HG22	2.20	0.41
1:E:56:SER:OG	1:E:58:TYR:CZ	2.72	0.41
1:F:12:LYS:HD3	1:F:58:TYR:CE1	2.56	0.41
1:F:48:LEU:HD22	1:F:48:LEU:HA	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:MET:HE3	1:A:128:VAL:HA	2.02	0.41
1:A:220:GLN:HB3	2:J:745:PHE:CE2	2.56	0.41
1:B:95:PRO:HG3	1:B:104:MET:CE	2.51	0.41
1:E:67:VAL:HG23	1:E:68:PHE:N	2.36	0.41
1:F:128:VAL:HG22	1:F:172:ILE:HD12	2.02	0.41
1:B:108:ASP:OD2	1:F:120:LYS:HE3	2.22	0.40
1:E:12:LYS:HD2	1:E:56:SER:OG	2.21	0.40
1:A:12:LYS:NZ	1:A:54:ASP:OD2	2.37	0.40
1:B:257:TRP:HZ2	1:D:37:GLU:HG2	1.87	0.40
1:C:103:ASN:OD1	1:C:105:VAL:HG13	2.21	0.40
1:C:206:CYS:HB3	1:C:241:LEU:HD23	2.04	0.40
1:A:59:ILE:HA	1:A:79:VAL:O	2.20	0.40
1:D:175:LEU:HD23	1:D:175:LEU:HA	1.81	0.40
1:E:126:LYS:HD3	1:E:126:LYS:N	2.36	0.40
1:F:61:ASP:HB3	1:F:62:PRO:HD3	2.04	0.40
1:A:190:ILE:CG2	1:A:191:ASN:N	2.84	0.40
1:A:208:THR:HG22	1:A:243:VAL:HG11	2.02	0.40
1:A:230:MET:C	1:A:232:GLN:N	2.79	0.40
1:B:261:CYS:C	1:B:284:THR:HG23	2.46	0.40
1:C:13:PHE:HA	1:C:59:ILE:HB	2.03	0.40
1:D:111:ILE:HG12	1:D:128:VAL:HG21	2.04	0.40
1:E:221:LEU:HD23	1:E:221:LEU:HA	1.84	0.40
1:F:26:ALA:HB1	1:F:84:VAL:CG2	2.52	0.40
1:A:28:GLU:HG3	1:A:29:SER:N	2.36	0.40
1:C:26:ALA:HB2	1:C:85:ILE:CD1	2.51	0.40
1:F:12:LYS:HE3	3:F:322:HOH:O	2.20	0.40
1:F:45:GLU:HG3	1:F:46:GLU:H	1.87	0.40
1:F:186:ARG:O	1:F:187:TYR:C	2.61	0.40

All (1) 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)
1:B:38:TYR:OH	3:A:301:HOH:O[2_554]	2.03	0.17

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	278/281 (99%)	268 (96%)	10 (4%)	0	100	100
1	B	274/281 (98%)	263 (96%)	11 (4%)	0	100	100
1	C	272/281 (97%)	258 (95%)	14 (5%)	0	100	100
1	D	276/281 (98%)	265 (96%)	11 (4%)	0	100	100
1	E	269/281 (96%)	257 (96%)	12 (4%)	0	100	100
1	F	271/281 (96%)	262 (97%)	9 (3%)	0	100	100
2	G	5/8 (62%)	5 (100%)	0	0	100	100
2	H	5/8 (62%)	5 (100%)	0	0	100	100
2	I	5/8 (62%)	5 (100%)	0	0	100	100
2	J	5/8 (62%)	4 (80%)	1 (20%)	0	100	100
All	All	1660/1718 (97%)	1592 (96%)	68 (4%)	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	257/259 (99%)	257 (100%)	0	100	100
1	B	246/259 (95%)	242 (98%)	4 (2%)	55	79
1	C	247/259 (95%)	244 (99%)	3 (1%)	63	82
1	D	251/259 (97%)	250 (100%)	1 (0%)	84	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	239/259 (92%)	237 (99%)	2 (1%)	73	86
1	F	247/259 (95%)	243 (98%)	4 (2%)	55	79
2	G	6/6 (100%)	6 (100%)	0	100	100
2	H	6/6 (100%)	6 (100%)	0	100	100
2	I	6/6 (100%)	5 (83%)	1 (17%)	2	12
2	J	6/6 (100%)	6 (100%)	0	100	100
All	All	1511/1578 (96%)	1496 (99%)	15 (1%)	68	84

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	50	ILE
1	B	94	VAL
1	B	181	GLU
1	B	183	LYS
1	C	21	LYS
1	C	23	PHE
1	C	181	GLU
1	D	178	LYS
1	E	53	ASN
1	E	131	MET
1	F	186	ARG
1	F	221	LEU
1	F	230	MET
1	F	283	LYS
2	I	750	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	260	HIS
1	B	91	GLN
1	B	258	ASN
1	B	260	HIS
1	B	277	GLN
1	C	265	GLN
1	D	53	ASN
1	D	82	GLN

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Mol	Chain	Res	Type
1	D	191	ASN
1	D	228	GLN
1	E	125	HIS
1	E	191	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SEP	J	749	2	8,9,10	1.86	2 (25%)	8,12,14	1.68	1 (12%)
2	SEP	G	749	2	8,9,10	0.76	0	8,12,14	1.90	1 (12%)
2	SEP	I	749	2	8,9,10	1.71	2 (25%)	8,12,14	1.82	2 (25%)
2	SEP	H	749	2	8,9,10	1.81	2 (25%)	8,12,14	1.84	3 (37%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	J	749	2	-	2/5/8/10	-
2	SEP	G	749	2	-	5/5/8/10	-
2	SEP	I	749	2	-	3/5/8/10	-
2	SEP	H	749	2	-	1/5/8/10	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	749	SEP	P-O1P	4.24	1.64	1.50
2	H	749	SEP	P-O1P	3.99	1.63	1.50
2	I	749	SEP	P-O1P	3.47	1.61	1.50
2	I	749	SEP	P-O3P	2.28	1.63	1.54
2	J	749	SEP	P-O2P	2.14	1.63	1.54
2	H	749	SEP	P-O3P	2.02	1.62	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	749	SEP	OG-CB-CA	4.50	112.53	108.14
2	J	749	SEP	OG-CB-CA	4.05	112.09	108.14
2	H	749	SEP	OG-CB-CA	3.21	111.27	108.14
2	I	749	SEP	P-OG-CB	-3.15	109.63	118.30
2	I	749	SEP	OG-CB-CA	2.81	110.88	108.14
2	H	749	SEP	O2P-P-OG	2.67	113.85	106.73
2	H	749	SEP	OG-P-O1P	2.11	112.41	106.47

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	749	SEP	N-CA-CB-OG
2	G	749	SEP	CA-CB-OG-P
2	G	749	SEP	CB-OG-P-O1P
2	G	749	SEP	CB-OG-P-O2P
2	G	749	SEP	CB-OG-P-O3P
2	J	749	SEP	N-CA-CB-OG
2	H	749	SEP	CA-CB-OG-P
2	J	749	SEP	CA-CB-OG-P
2	I	749	SEP	CB-OG-P-O1P
2	I	749	SEP	CB-OG-P-O2P
2	I	749	SEP	CA-CB-OG-P

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

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	280/281 (99%)	-0.00	1 (0%) 88 76	49, 69, 103, 122	0
1	B	278/281 (98%)	-0.01	4 (1%) 73 51	49, 67, 100, 116	0
1	C	276/281 (98%)	0.02	6 (2%) 62 39	50, 69, 98, 116	0
1	D	280/281 (99%)	0.01	0 100 100	54, 74, 101, 122	0
1	E	273/281 (97%)	0.25	11 (4%) 42 23	50, 81, 120, 136	0
1	F	275/281 (97%)	0.30	7 (2%) 58 35	54, 84, 117, 135	0
2	G	7/8 (87%)	0.06	0 100 100	50, 56, 76, 85	0
2	H	7/8 (87%)	0.11	0 100 100	59, 64, 81, 89	0
2	I	7/8 (87%)	-0.08	0 100 100	49, 59, 69, 79	0
2	J	7/8 (87%)	0.46	1 (14%) 6 3	61, 62, 78, 94	0
All	All	1690/1718 (98%)	0.09	30 (1%) 67 44	49, 74, 110, 136	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	60	CYS	4.1
1	F	284	THR	3.8
1	B	52	GLU	3.7
1	E	186	ARG	3.7
1	F	16	SER	3.6
1	F	256	ARG	3.6
1	E	56	SER	3.5
1	F	252	GLU	3.2
1	C	19	ASN	3.2
1	E	67	VAL	2.9
1	C	15	LYS	2.9
1	C	275	PHE	2.8
1	F	82	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	48	LEU	2.8
1	B	286	PRO	2.7
1	F	139	LEU	2.7
1	C	35	SER	2.5
1	F	20	SER	2.5
1	C	22	CYS	2.5
1	B	51	LYS	2.5
1	A	137	ARG	2.5
1	B	61	ASP	2.4
1	E	55	ARG	2.4
2	J	743	SER	2.3
1	E	59	ILE	2.3
1	E	42	ILE	2.2
1	E	66	VAL	2.2
1	E	139	LEU	2.1
1	E	52	GLU	2.1
1	E	62	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	SEP	G	749	10/11	0.68	0.17	59,64,75,81	0
2	SEP	H	749	10/11	0.92	0.09	61,75,90,92	0
2	SEP	J	749	10/11	0.93	0.08	59,63,67,73	0
2	SEP	I	749	10/11	0.97	0.05	54,64,71,87	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.