



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:25 PM GMT

PDB ID : 1K8G
Title : Crystal Structure of the N-terminal domain of Oxytricha nova telomere end binding protein alpha subunit both uncomplexed and complexed with telomeric ssDNA
Authors : Classen, S.; Ruggles, J.A.; Schultz, S.C.
Deposited on : 2001-10-24
Resolution : 2.60 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

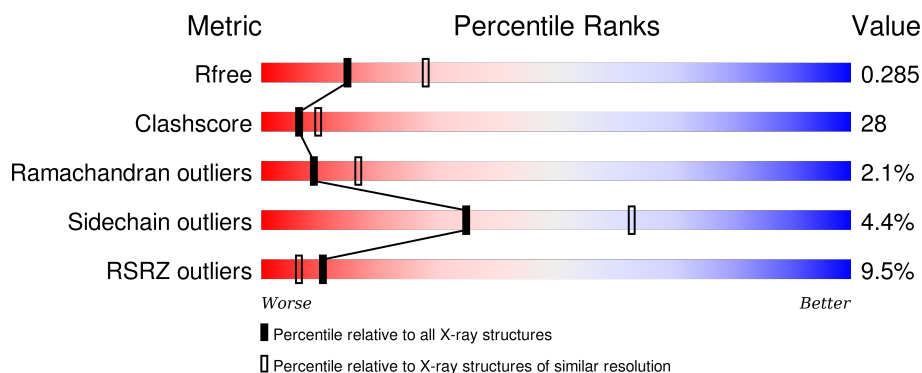
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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. 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	D	6	<div> <div>17%</div> <div>67%</div> <div>33%</div> </div>
1	E	6	<div> <div>17%</div> <div>17%</div> <div>33%</div> <div>17%</div> <div>33%</div> </div>
2	A	320	<div> <div>8%</div> <div>47%</div> <div>34%</div> <div>•</div> <div>16%</div> </div>
2	B	320	<div> <div>9%</div> <div>41%</div> <div>42%</div> <div>•</div> <div>14%</div> </div>
2	C	320	<div> <div>7%</div> <div>43%</div> <div>42%</div> <div>•</div> <div>13%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7040 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.

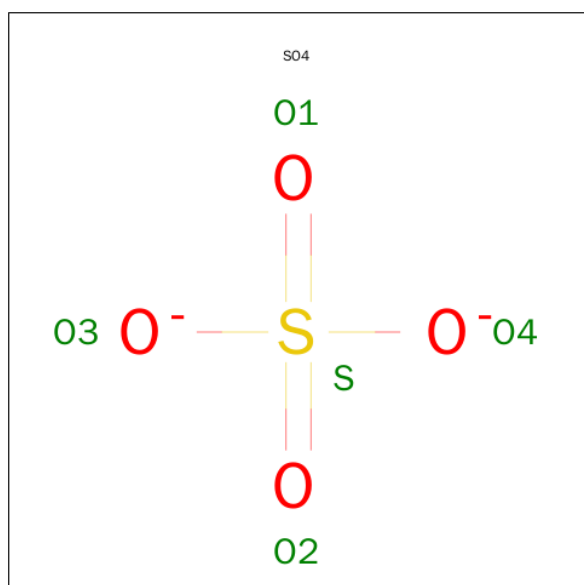
- Molecule 1 is a DNA chain called 5'-D(TP*TP*GP*GP*GP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	4	Total	C	N	O	P	0	0	0
			85	40	20	22	3			
1	E	4	Total	C	N	O	P	0	0	0
			85	40	20	22	3			

- Molecule 2 is a protein called Telomere-Binding Protein alpha Subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	269	Total	C	N	O	S	0	0	0
			2177	1395	367	413	2			
2	B	275	Total	C	N	O	S	0	0	0
			2231	1424	379	426	2			
2	C	278	Total	C	N	O	S	0	0	0
			2249	1435	383	429	2			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0

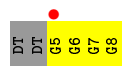
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	59	Total O 59 59	0	0
4	B	55	Total O 55 55	0	0
4	C	66	Total O 66 66	0	0
4	D	2	Total O 2 2	0	0
4	E	1	Total O 1 1	0	0

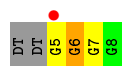
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

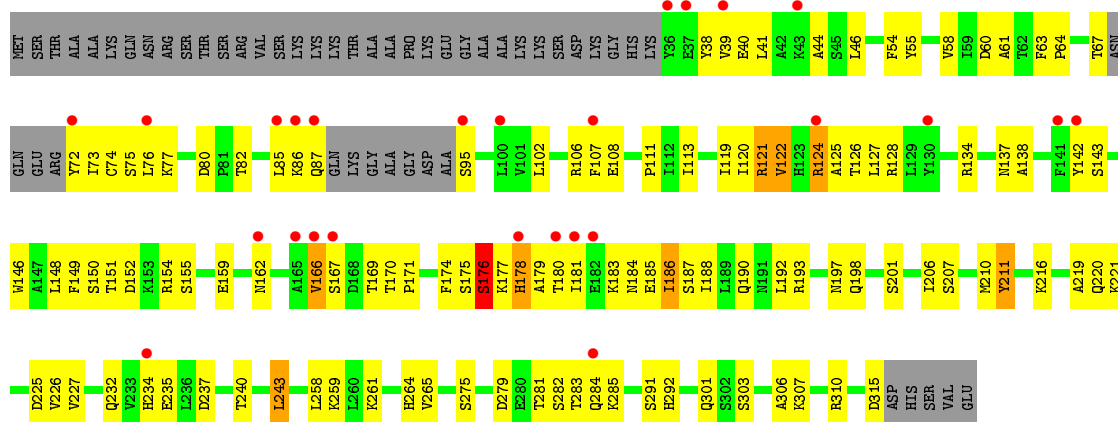
- Molecule 1: 5'-D(TP*TP*GP*GP*GP*G)-3'



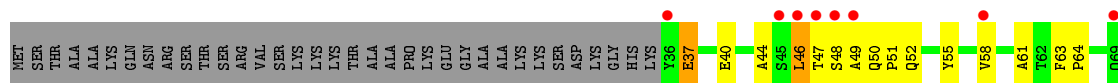
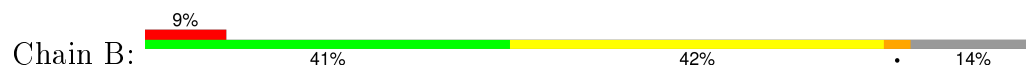
- Molecule 1: 5'-D(TP*TP*GP*GP*GP*G)-3'

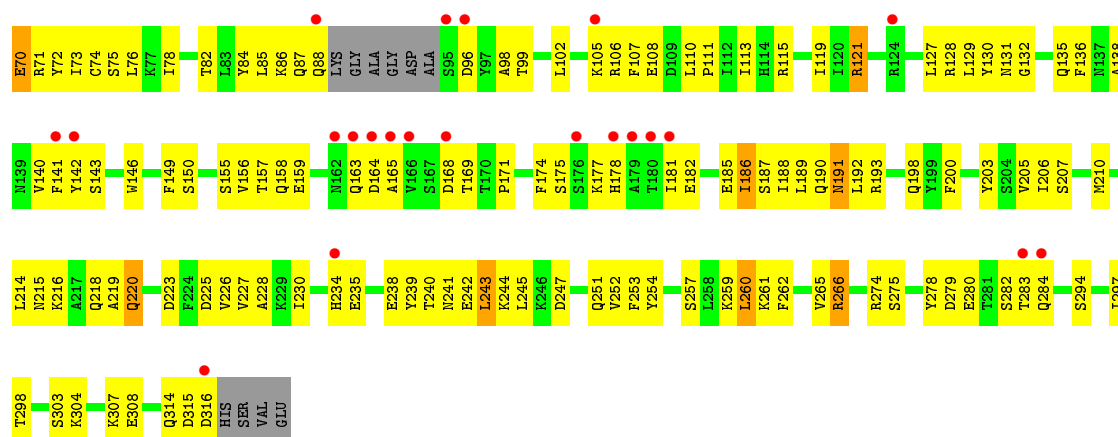


- Molecule 2: Telomere-Binding Protein alpha Subunit

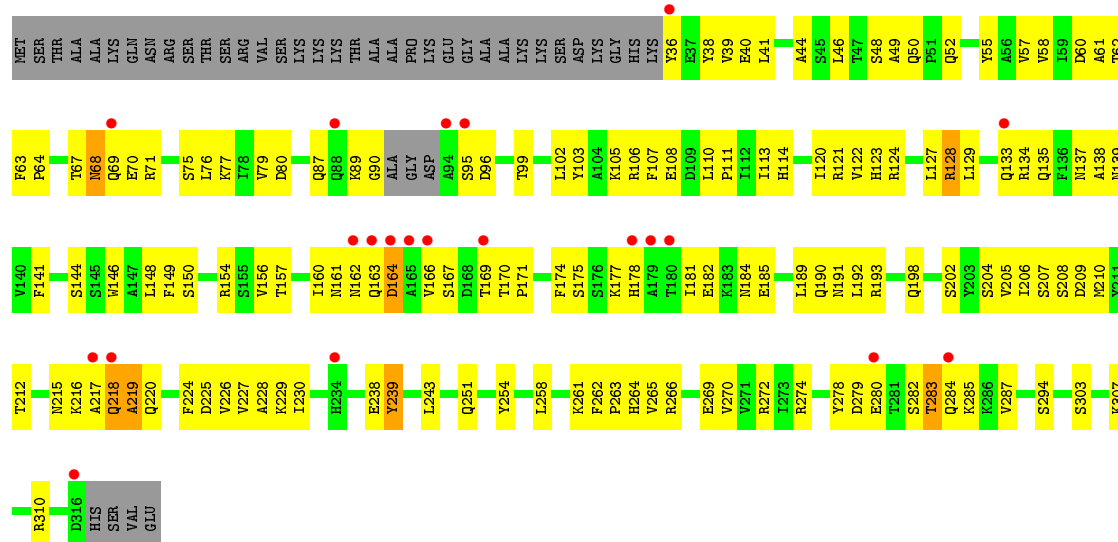
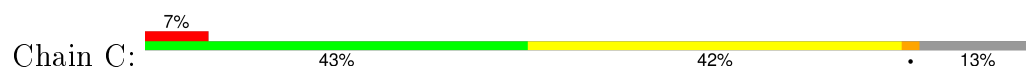


- Molecule 2: Telomere-Binding Protein alpha Subunit





• Molecule 2: Telomere-Binding Protein alpha Subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.18Å 121.18Å 137.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.60 29.96 – 2.61	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.60) 99.8 (29.96-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.285 0.245 , 0.285	Depositor DCC
R_{free} test set	3583 reflections (9.96%)	DCC
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.159	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.8	EDS
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 35965 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7040	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.62% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: SO4

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	0.51	0/96	0.77	0/148
1	E	0.94	0/96	1.31	3/148 (2.0%)
2	A	0.47	0/2221	0.76	0/3008
2	B	0.48	0/2276	0.73	0/3083
2	C	0.47	0/2294	0.74	0/3106
All	All	0.48	0/6983	0.75	3/9493 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	E	6	DG	N9-C4-C5	8.22	108.69	105.40
1	E	6	DG	C8-N9-C4	-6.90	103.64	106.40
1	E	6	DG	C4-C5-N7	-5.01	108.80	110.80

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	D	85	0	46	7	0
1	E	85	0	46	6	0
2	A	2177	0	2164	123	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2231	0	2210	135	0
2	C	2249	0	2231	122	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
4	A	59	0	0	1	0
4	B	55	0	0	6	0
4	C	66	0	0	3	0
4	D	2	0	0	0	0
4	E	1	0	0	0	0
All	All	7040	0	6697	376	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 28.

All (376) 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:C:169:THR:HG21	2:C:181:ILE:HG12	1.29	1.08
2:A:85:LEU:HD13	2:A:95:SER:HA	1.42	1.02
2:B:141:PHE:HD2	2:B:142:TYR:HD2	1.09	1.01
2:A:169:THR:HG23	2:A:181:ILE:HD11	1.45	0.98
2:A:206:ILE:HA	2:A:210:MET:CE	1.94	0.98
2:C:154:ARG:NH1	2:C:162:ASN:HA	1.80	0.96
2:B:87:GLN:O	2:B:88:GLN:HG2	1.66	0.94
2:C:58:VAL:HG11	2:C:113:ILE:HD13	1.48	0.94
2:C:283:THR:HG23	2:C:284:GLN:H	1.38	0.88
2:B:141:PHE:HD2	2:B:142:TYR:CD2	1.91	0.87
2:A:169:THR:CG2	2:A:181:ILE:HD11	2.03	0.87
2:A:279:ASP:OD1	2:A:281:THR:HG22	1.75	0.86
2:C:154:ARG:HH12	2:C:162:ASN:HA	1.36	0.86
2:C:181:ILE:HD12	2:C:185:GLU:OE1	1.79	0.83
2:A:176:SER:CB	2:A:179:ALA:HB2	2.07	0.83
2:C:218:GLN:HG3	2:C:278:TYR:CE2	2.14	0.82
2:B:71:ARG:NH1	2:B:105:LYS:HE2	1.94	0.82
2:A:176:SER:OG	2:A:179:ALA:HB2	1.80	0.82
2:A:121:ARG:HH12	2:A:181:ILE:CG2	1.93	0.82
2:B:141:PHE:CD2	2:B:142:TYR:HD2	1.97	0.82
2:C:105:LYS:HE3	2:C:141:PHE:HB2	1.63	0.80
2:B:85:LEU:HD13	2:B:96:ASP:OD2	1.81	0.80
2:B:158:GLN:HG2	2:B:163:GLN:HE21	1.46	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:121:ARG:HH12	2:A:181:ILE:HG22	1.47	0.80
2:A:206:ILE:HA	2:A:210:MET:HE3	1.63	0.80
2:A:124:ARG:HD3	2:A:143:SER:O	1.83	0.79
2:A:180:THR:HG22	2:A:181:ILE:N	1.96	0.79
2:B:70:GLU:O	2:B:71:ARG:HD3	1.82	0.79
1:E:6:DG:H3'	1:E:7:DG:C5'	2.13	0.79
2:C:102:LEU:HD23	2:C:138:ALA:HB3	1.65	0.78
2:B:230:ILE:HD13	2:B:265:VAL:HG13	1.66	0.78
2:A:180:THR:HG22	2:A:181:ILE:H	1.48	0.78
2:B:61:ALA:CB	2:B:76:LEU:HD22	2.14	0.78
2:C:230:ILE:HD13	2:C:265:VAL:HG13	1.68	0.74
2:A:181:ILE:HB	2:A:185:GLU:OE1	1.87	0.73
2:C:169:THR:CG2	2:C:181:ILE:HG12	2.14	0.73
2:C:67:THR:OG1	2:C:71:ARG:HB3	1.89	0.73
2:B:61:ALA:HB2	2:B:76:LEU:HD22	1.69	0.72
2:B:266:ARG:HG2	2:B:266:ARG:HH11	1.54	0.72
2:A:75:SER:C	2:A:76:LEU:HD23	2.10	0.71
2:A:197:ASN:O	2:A:201:SER:HB3	1.90	0.71
2:A:154:ARG:NH1	2:A:162:ASN:HA	2.05	0.71
1:D:5:DG:N3	1:D:5:DG:H2'	2.06	0.70
2:C:206:ILE:HA	2:C:210:MET:CE	2.21	0.70
2:C:128:ARG:NH1	2:C:135:GLN:OE1	2.23	0.70
2:A:225:ASP:H	2:B:87:GLN:HE22	1.38	0.70
1:D:6:DG:H3'	1:D:7:DG:C5'	2.21	0.70
2:C:61:ALA:CB	2:C:76:LEU:HD22	2.22	0.70
2:B:216:LYS:HG2	2:C:202:SER:HB2	1.74	0.69
2:A:243:LEU:HD11	2:A:265:VAL:HG11	1.74	0.69
2:B:169:THR:O	2:B:169:THR:HG22	1.92	0.69
2:B:71:ARG:HH12	2:B:105:LYS:HE2	1.57	0.68
2:C:111:PRO:HG3	2:C:146:TRP:CD2	2.28	0.68
2:A:85:LEU:HB2	2:A:95:SER:HB2	1.76	0.68
2:A:154:ARG:HH12	2:A:162:ASN:HA	1.59	0.68
2:A:85:LEU:HB2	2:A:95:SER:CB	2.23	0.68
2:C:160:ILE:HG12	2:C:266:ARG:HH21	1.59	0.67
2:B:304:LYS:HE3	2:B:308:GLU:OE2	1.95	0.67
2:B:130:TYR:O	2:B:131:ASN:OD1	2.12	0.67
2:C:189:LEU:HD21	2:C:193:ARG:NH2	2.10	0.66
2:B:206:ILE:HA	2:B:210:MET:HE3	1.75	0.66
2:C:279:ASP:HB3	2:C:282:SER:HB2	1.75	0.66
2:B:243:LEU:HD11	2:B:265:VAL:HG11	1.77	0.66
2:B:207:SER:H	2:B:210:MET:HE2	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:105:LYS:HD2	2:B:141:PHE:CE1	2.31	0.65
2:A:87:GLN:HE22	2:C:225:ASP:H	1.44	0.65
2:B:70:GLU:C	2:B:71:ARG:HD3	2.17	0.65
1:D:6:DG:H3'	1:D:7:DG:H5"	1.79	0.64
2:C:156:VAL:HG13	2:C:157:THR:N	2.12	0.64
2:B:206:ILE:HA	2:B:210:MET:CE	2.27	0.64
2:B:257:SER:HB2	2:B:262:PHE:HD1	1.62	0.64
2:C:207:SER:HB3	2:C:210:MET:HG3	1.80	0.64
2:C:166:VAL:HB	2:C:170:THR:HG21	1.79	0.64
2:B:99:THR:HG21	4:B:343:HOH:O	1.97	0.64
2:B:191:ASN:N	2:B:191:ASN:HD22	1.95	0.63
2:B:155:SER:HB3	2:B:158:GLN:HG3	1.80	0.63
2:C:61:ALA:HB1	2:C:76:LEU:HD22	1.80	0.63
2:A:206:ILE:HA	2:A:210:MET:HE2	1.80	0.63
2:A:102:LEU:HD23	2:A:138:ALA:HB3	1.81	0.63
2:A:149:PHE:CZ	2:A:171:PRO:HG3	2.34	0.63
2:A:180:THR:CG2	2:A:181:ILE:H	2.11	0.62
2:B:72:TYR:CE2	2:B:107:PHE:HB2	2.34	0.62
2:C:160:ILE:HG23	2:C:266:ARG:NH2	2.15	0.62
2:C:111:PRO:HG3	2:C:146:TRP:CE2	2.34	0.62
2:B:168:ASP:O	2:B:193:ARG:NH2	2.31	0.62
2:B:177:LYS:O	2:B:178:HIS:HD2	1.84	0.61
2:C:204:SER:O	2:C:206:ILE:N	2.34	0.60
2:C:238:GLU:HB2	2:C:239:TYR:CE1	2.36	0.60
2:A:185:GLU:O	2:A:186:ILE:C	2.40	0.60
2:A:180:THR:CG2	2:A:181:ILE:N	2.64	0.60
2:C:206:ILE:HA	2:C:210:MET:HE2	1.83	0.60
2:A:221:LYS:HG2	4:B:340:HOH:O	2.01	0.60
2:A:142:TYR:HD2	2:A:143:SER:HG	1.50	0.60
2:A:282:SER:HB2	2:A:285:LYS:O	2.01	0.59
2:B:37:GLU:O	2:B:37:GLU:HG2	2.00	0.59
2:C:226:VAL:HG12	2:C:227:VAL:N	2.17	0.59
2:A:177:LYS:O	2:A:178:HIS:HB2	2.02	0.59
2:B:40:GLU:OE1	2:B:82:THR:HG21	2.02	0.59
2:B:61:ALA:HB1	2:B:76:LEU:HD22	1.84	0.59
2:C:215:ASN:OD1	2:C:216:LYS:HG3	2.02	0.59
2:B:78:ILE:HG13	2:B:98:ALA:HB3	1.85	0.59
2:B:111:PRO:HG3	2:B:146:TRP:CD2	2.38	0.59
2:A:186:ILE:O	2:A:190:GLN:HG2	2.03	0.59
1:E:6:DG:H3'	1:E:7:DG:H5"	1.84	0.59
2:C:217:ALA:HB1	2:C:224:PHE:CZ	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:226:VAL:HG12	2:A:227:VAL:N	2.18	0.58
2:A:206:ILE:HD13	2:A:210:MET:HE1	1.84	0.58
2:A:150:SER:O	2:A:193:ARG:NH1	2.36	0.58
2:C:70:GLU:O	2:C:105:LYS:HA	2.03	0.58
2:C:243:LEU:HD11	2:C:265:VAL:HG11	1.85	0.58
2:B:146:TRP:CE2	2:B:175:SER:HB3	2.39	0.57
2:C:212:THR:HG21	2:C:220:GLN:HE21	1.68	0.57
2:C:64:PRO:HB2	2:C:107:PHE:CE1	2.40	0.57
2:A:185:GLU:O	2:A:188:ILE:N	2.37	0.57
1:D:6:DG:N2	2:A:128:ARG:NH2	2.52	0.57
2:C:206:ILE:HA	2:C:210:MET:HE1	1.87	0.57
2:A:146:TRP:NE1	2:A:175:SER:HB3	2.20	0.57
2:C:129:LEU:HD12	2:C:133:GLN:O	2.05	0.57
2:A:74:CYS:SG	2:A:76:LEU:HD21	2.45	0.57
2:C:190:GLN:OE1	2:C:190:GLN:HA	2.04	0.57
2:B:235:GLU:HG3	4:B:327:HOH:O	2.04	0.57
2:A:177:LYS:O	2:A:178:HIS:CB	2.53	0.56
2:A:225:ASP:H	2:B:87:GLN:NE2	2.03	0.56
2:A:185:GLU:O	2:A:187:SER:N	2.39	0.56
2:A:60:ASP:HB3	2:A:77:LYS:HB2	1.88	0.56
2:A:85:LEU:HB2	2:A:95:SER:HA	1.88	0.56
2:A:85:LEU:CD1	2:A:95:SER:HA	2.25	0.56
2:B:46:LEU:HD22	2:B:127:LEU:CD2	2.36	0.56
2:B:216:LYS:O	2:B:220:GLN:OE1	2.24	0.56
2:A:167:SER:O	2:A:170:THR:HB	2.07	0.55
2:B:274:ARG:O	2:B:275:SER:HB3	2.06	0.55
2:C:114:HIS:HB2	2:C:156:VAL:HG11	1.87	0.55
2:C:44:ALA:HB1	2:C:52:GLN:OE1	2.06	0.55
2:C:61:ALA:HB2	2:C:76:LEU:HD22	1.88	0.55
2:B:102:LEU:HD23	2:B:138:ALA:HB3	1.87	0.55
2:B:141:PHE:CD2	2:B:142:TYR:CD2	2.82	0.55
2:C:177:LYS:O	2:C:178:HIS:CG	2.60	0.54
2:C:48:SER:C	2:C:50:GLN:H	2.11	0.54
2:B:181:ILE:O	2:B:181:ILE:HG13	2.07	0.54
2:C:207:SER:HB3	2:C:210:MET:CG	2.38	0.54
2:C:239:TYR:O	2:C:258:LEU:HD12	2.08	0.54
2:A:72:TYR:HE2	2:A:107:PHE:CE1	2.25	0.54
2:C:156:VAL:HG13	2:C:157:THR:H	1.73	0.53
2:A:58:VAL:HG11	2:A:113:ILE:HG12	1.90	0.53
2:C:149:PHE:CE2	2:C:169:THR:O	2.60	0.53
2:B:280:GLU:HA	2:B:280:GLU:OE1	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:121:ARG:HH12	2:A:181:ILE:HG21	1.74	0.53
2:C:99:THR:HG21	4:C:373:HOH:O	2.09	0.53
2:A:197:ASN:O	2:A:201:SER:CB	2.57	0.53
2:B:70:GLU:O	2:B:105:LYS:HA	2.09	0.53
2:B:155:SER:O	2:B:159:GLU:HG2	2.09	0.53
2:B:234:HIS:HB3	2:B:242:GLU:HB3	1.89	0.53
2:C:167:SER:O	2:C:170:THR:HG22	2.09	0.53
2:B:297:ILE:HG22	2:B:298:THR:O	2.09	0.52
2:C:60:ASP:HB3	2:C:77:LYS:HB2	1.91	0.52
2:B:218:GLN:HB2	4:B:351:HOH:O	2.08	0.52
2:A:63:PHE:CD2	2:A:64:PRO:HD2	2.44	0.52
2:B:186:ILE:O	2:B:190:GLN:HB2	2.09	0.52
2:B:129:LEU:HD21	2:B:132:GLY:CA	2.40	0.52
2:B:73:ILE:HG22	2:B:74:CYS:N	2.23	0.52
2:A:152:ASP:OD2	2:A:301:GLN:HG3	2.09	0.52
2:A:207:SER:HB3	2:A:210:MET:HG3	1.92	0.52
2:B:149:PHE:CZ	2:B:171:PRO:HG3	2.44	0.52
2:B:87:GLN:O	2:B:88:GLN:CG	2.50	0.52
2:B:129:LEU:HD21	2:B:132:GLY:HA2	1.92	0.52
2:A:126:THR:CG2	2:A:137:ASN:HB2	2.40	0.52
2:C:218:GLN:O	2:C:220:GLN:N	2.42	0.51
2:C:160:ILE:HG23	2:C:266:ARG:HH22	1.75	0.51
2:A:176:SER:HB2	2:A:179:ALA:HB2	1.89	0.51
2:B:266:ARG:HG2	2:B:266:ARG:NH1	2.22	0.51
2:C:103:TYR:HE1	2:C:137:ASN:HD22	1.57	0.51
2:A:226:VAL:CG1	2:A:227:VAL:N	2.73	0.51
2:B:140:VAL:HG12	2:B:140:VAL:O	2.11	0.51
2:C:124:ARG:NH1	2:C:124:ARG:HG3	2.25	0.51
2:C:254:TYR:CE1	2:C:285:LYS:HE3	2.45	0.51
2:C:124:ARG:HH11	2:C:124:ARG:HG3	1.75	0.51
2:B:219:ALA:HB2	2:C:198:GLN:HB3	1.92	0.51
2:A:85:LEU:HB2	2:A:95:SER:CA	2.41	0.50
2:B:261:LYS:HE2	2:B:262:PHE:CZ	2.46	0.50
2:A:166:VAL:HG12	2:A:167:SER:N	2.25	0.50
2:A:206:ILE:CA	2:A:210:MET:HE3	2.39	0.50
2:C:68:ASN:ND2	2:C:71:ARG:HB2	2.27	0.50
2:A:174:PHE:CE1	2:A:178:HIS:HA	2.46	0.50
2:B:215:ASN:HA	2:B:251:GLN:HG3	1.94	0.50
2:B:230:ILE:CD1	2:B:265:VAL:HG13	2.40	0.50
2:B:207:SER:H	2:B:210:MET:CE	2.24	0.50
2:B:259:LYS:HD2	4:B:327:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:228:ALA:O	2:C:270:VAL:HA	2.10	0.50
2:B:44:ALA:HA	2:B:52:GLN:HE22	1.77	0.50
1:D:8:DG:O6	2:A:275:SER:HB2	2.11	0.50
2:B:218:GLN:HG3	2:B:278:TYR:CE1	2.47	0.50
2:A:146:TRP:CE2	2:A:175:SER:HB3	2.47	0.49
2:B:58:VAL:HG11	2:B:113:ILE:HG12	1.94	0.49
2:C:307:LYS:HG2	2:C:310:ARG:NH2	2.26	0.49
2:C:239:TYR:CD1	2:C:239:TYR:N	2.79	0.49
1:E:5:DG:H4'	2:B:73:ILE:HD11	1.93	0.49
2:B:257:SER:HA	4:B:328:HOH:O	2.12	0.49
2:C:36:TYR:HB3	2:C:38:TYR:CE1	2.47	0.49
2:C:181:ILE:HG23	2:C:181:ILE:O	2.12	0.49
2:B:228:ALA:HB2	2:B:247:ASP:HB3	1.94	0.49
2:A:75:SER:O	2:A:76:LEU:HD23	2.12	0.49
2:B:164:ASP:OD1	2:B:165:ALA:N	2.46	0.49
2:B:46:LEU:HD22	2:B:127:LEU:HD23	1.93	0.49
2:A:67:THR:HG21	2:A:73:ILE:HD12	1.95	0.49
2:A:61:ALA:CB	2:A:76:LEU:HD22	2.43	0.49
2:B:189:LEU:HD21	2:B:193:ARG:NH2	2.28	0.48
2:C:123:HIS:HB3	2:C:124:ARG:NH1	2.28	0.48
2:C:282:SER:OG	2:C:285:LYS:O	2.30	0.48
2:A:119:ILE:HD13	2:A:192:LEU:HD23	1.94	0.48
2:C:68:ASN:OD1	2:C:68:ASN:C	2.52	0.48
2:B:64:PRO:HG3	2:B:111:PRO:O	2.13	0.48
2:B:119:ILE:CG2	2:B:192:LEU:HD23	2.44	0.48
2:B:223:ASP:O	2:C:89:LYS:NZ	2.45	0.48
2:B:158:GLN:HG2	2:B:163:GLN:NE2	2.21	0.48
2:B:146:TRP:NE1	2:B:175:SER:HB3	2.28	0.48
2:C:46:LEU:HD23	2:C:127:LEU:HG	1.95	0.48
2:C:184:ASN:OD1	2:C:185:GLU:N	2.46	0.48
2:C:303:SER:O	2:C:307:LYS:HG3	2.14	0.48
2:C:264:HIS:O	2:C:266:ARG:HG2	2.12	0.48
2:A:119:ILE:CG2	2:A:120:ILE:N	2.77	0.48
2:C:62:THR:HG22	2:C:274:ARG:NH1	2.29	0.48
2:C:229:LYS:HA	2:C:269:GLU:O	2.14	0.48
2:B:283:THR:HG23	2:B:283:THR:O	2.14	0.47
2:A:169:THR:O	2:A:169:THR:HG22	2.14	0.47
2:B:207:SER:HB2	2:B:210:MET:HG3	1.96	0.47
2:B:63:PHE:CD2	2:B:261:LYS:HE3	2.50	0.47
2:B:168:ASP:N	2:B:168:ASP:OD1	2.36	0.47
2:C:106:ARG:CZ	2:C:108:GLU:OE2	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:206:ILE:HD13	2:C:210:MET:CE	2.45	0.47
2:C:48:SER:O	2:C:50:GLN:N	2.42	0.47
2:B:279:ASP:O	2:B:282:SER:OG	2.23	0.47
2:B:105:LYS:CD	2:B:141:PHE:CE1	2.98	0.47
2:A:225:ASP:N	2:B:87:GLN:HE22	2.08	0.47
2:C:283:THR:HG23	2:C:284:GLN:N	2.17	0.47
2:A:306:ALA:O	2:A:310:ARG:HG3	2.14	0.47
2:A:73:ILE:CG2	2:A:74:CYS:N	2.77	0.47
2:C:265:VAL:O	2:C:266:ARG:HD3	2.14	0.47
2:C:75:SER:O	2:C:76:LEU:HD23	2.13	0.47
2:B:185:GLU:O	2:B:188:ILE:N	2.48	0.47
2:C:39:VAL:HG23	4:C:343:HOH:O	2.15	0.47
2:C:63:PHE:HB2	2:C:294:SER:O	2.15	0.47
2:A:124:ARG:HD3	2:A:143:SER:C	2.33	0.47
2:B:191:ASN:ND2	2:B:191:ASN:N	2.62	0.47
2:B:200:PHE:CZ	2:B:205:VAL:HG11	2.50	0.47
2:A:119:ILE:HG22	2:A:120:ILE:N	2.29	0.47
2:C:207:SER:O	2:C:209:ASP:N	2.49	0.46
2:B:55:TYR:HB3	2:B:192:LEU:HD22	1.97	0.46
1:E:6:DG:H3'	1:E:7:DG:H5'	1.97	0.46
2:A:206:ILE:HD13	2:A:210:MET:CE	2.46	0.46
2:B:218:GLN:HG3	2:B:278:TYR:CZ	2.51	0.46
2:C:39:VAL:HG12	2:C:40:GLU:O	2.16	0.46
2:A:142:TYR:CD2	2:A:143:SER:N	2.83	0.46
2:C:261:LYS:C	2:C:263:PRO:HD3	2.35	0.46
2:B:314:GLN:OE1	2:B:314:GLN:HA	2.16	0.46
2:C:164:ASP:N	2:C:164:ASP:OD1	2.49	0.46
2:B:315:ASP:OD1	2:B:316:ASP:N	2.49	0.46
2:A:198:GLN:HB3	2:C:219:ALA:HB2	1.97	0.46
2:B:127:LEU:HD13	2:B:136:PHE:CE2	2.51	0.46
2:A:183:LYS:C	2:A:185:GLU:N	2.69	0.46
2:A:61:ALA:HB1	2:A:76:LEU:HD22	1.98	0.46
2:A:41:LEU:O	2:A:134:ARG:NH1	2.48	0.46
2:B:64:PRO:HB3	2:B:110:LEU:HB2	1.98	0.45
2:A:219:ALA:HB2	2:B:198:GLN:HB3	1.98	0.45
2:B:284:GLN:HG2	2:B:284:GLN:O	2.16	0.45
2:A:216:LYS:O	2:A:220:GLN:HG2	2.16	0.45
2:A:207:SER:H	2:A:210:MET:HE2	1.82	0.45
2:C:46:LEU:HD23	2:C:127:LEU:CD2	2.46	0.45
2:C:146:TRP:CE2	2:C:175:SER:HB3	2.52	0.45
2:A:46:LEU:HD22	2:A:127:LEU:CD2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:235:GLU:OE2	2:A:259:LYS:NZ	2.50	0.45
2:C:218:GLN:HG3	2:C:278:TYR:CD2	2.50	0.45
2:C:182:GLU:HA	2:C:182:GLU:OE1	2.17	0.45
2:A:95:SER:N	2:A:207:SER:HB2	2.32	0.45
2:C:89:LYS:HB2	2:C:90:GLY:H	1.66	0.45
2:A:39:VAL:HG11	2:A:44:ALA:HA	1.98	0.45
2:A:206:ILE:CD1	2:A:210:MET:HE1	2.46	0.45
2:B:73:ILE:CG2	2:B:74:CYS:N	2.80	0.45
2:C:156:VAL:CG1	2:C:157:THR:N	2.79	0.45
2:C:139:ASN:HB3	2:C:141:PHE:CZ	2.52	0.44
2:B:72:TYR:CE1	2:B:107:PHE:N	2.86	0.44
2:A:55:TYR:HB3	2:A:192:LEU:HD22	2.00	0.44
2:B:182:GLU:OE1	2:B:182:GLU:HA	2.17	0.44
2:A:121:ARG:NH1	2:A:181:ILE:HG22	2.23	0.44
1:E:6:DG:H5'	2:B:73:ILE:HD13	1.99	0.44
2:C:146:TRP:NE1	2:C:175:SER:HB3	2.32	0.44
2:C:226:VAL:CG1	2:C:227:VAL:N	2.79	0.44
2:B:235:GLU:HA	2:B:241:ASN:ND2	2.33	0.44
2:B:61:ALA:HB1	2:B:76:LEU:CD2	2.46	0.44
2:C:75:SER:C	2:C:76:LEU:HD23	2.38	0.44
2:B:99:THR:HG22	2:B:135:GLN:HG3	1.98	0.44
2:B:185:GLU:O	2:B:187:SER:N	2.51	0.44
2:C:55:TYR:HB3	2:C:192:LEU:HD22	1.99	0.44
2:A:149:PHE:CE2	2:A:171:PRO:HG3	2.53	0.44
2:C:251:GLN:HA	2:C:251:GLN:OE1	2.18	0.44
2:B:244:LYS:HG3	2:B:254:TYR:CE2	2.52	0.44
2:A:183:LYS:O	2:A:185:GLU:N	2.51	0.44
2:C:280:GLU:N	2:C:280:GLU:OE1	2.50	0.44
2:B:225:ASP:H	2:C:87:GLN:HE22	1.65	0.44
1:D:7:DG:H1	2:A:225:ASP:CG	2.21	0.44
2:A:120:ILE:HG13	2:A:148:LEU:HD23	2.00	0.44
2:B:226:VAL:HG12	2:B:227:VAL:N	2.33	0.44
2:C:150:SER:O	2:C:193:ARG:NH1	2.51	0.43
2:C:285:LYS:O	2:C:287:VAL:HG23	2.18	0.43
2:C:62:THR:CG2	2:C:274:ARG:NH1	2.81	0.43
2:C:262:PHE:N	2:C:263:PRO:HD3	2.32	0.43
2:B:128:ARG:HG2	2:B:129:LEU:N	2.33	0.43
2:A:283:THR:OG1	2:A:284:GLN:N	2.49	0.43
2:B:105:LYS:HG3	2:B:141:PHE:CD1	2.54	0.43
2:B:214:LEU:HD13	2:B:253:PHE:CD1	2.54	0.43
2:C:61:ALA:HB1	2:C:76:LEU:CD2	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:46:LEU:CD2	2:B:127:LEU:HD21	2.49	0.43
2:B:129:LEU:HD23	2:B:132:GLY:H	1.84	0.43
2:A:211:TYR:N	2:A:211:TYR:CD2	2.87	0.43
2:A:38:TYR:CZ	2:A:185:GLU:OE2	2.72	0.43
2:B:260:LEU:HD12	2:B:260:LEU:HA	1.71	0.42
2:A:76:LEU:HD23	2:A:76:LEU:N	2.34	0.42
2:B:185:GLU:O	2:B:186:ILE:C	2.58	0.42
2:A:216:LYS:HD3	2:B:203:TYR:CZ	2.55	0.42
2:B:48:SER:O	2:B:50:GLN:N	2.52	0.42
2:B:50:GLN:HA	2:B:51:PRO:HD3	1.77	0.42
2:B:146:TRP:CZ3	2:B:174:PHE:HA	2.54	0.42
2:B:129:LEU:CD2	2:B:132:GLY:H	2.32	0.42
2:B:119:ILE:HG23	2:B:192:LEU:HD23	2.02	0.42
2:C:39:VAL:HG12	2:C:40:GLU:N	2.35	0.42
1:E:6:DG:C3'	1:E:7:DG:C5'	2.91	0.42
2:C:207:SER:C	2:C:209:ASP:H	2.23	0.42
2:C:254:TYR:CZ	2:C:285:LYS:HE3	2.54	0.42
2:B:102:LEU:CD2	2:B:138:ALA:HB3	2.50	0.42
2:A:80:ASP:OD1	2:A:80:ASP:C	2.58	0.42
2:B:115:ARG:HB2	2:B:156:VAL:HG21	2.02	0.42
2:A:151:THR:HA	2:A:193:ARG:HD3	2.02	0.42
2:A:54:PHE:O	2:A:121:ARG:HA	2.20	0.42
2:A:169:THR:HG22	2:A:181:ILE:HD11	1.93	0.42
2:A:72:TYR:N	2:A:72:TYR:CD1	2.87	0.42
2:B:297:ILE:HG22	2:B:298:THR:N	2.34	0.42
2:B:245:LEU:O	2:B:252:VAL:HA	2.20	0.42
2:A:86:LYS:O	2:A:87:GLN:C	2.58	0.41
2:A:170:THR:HA	2:A:171:PRO:HD3	1.79	0.41
2:C:161:ASN:HB2	2:C:163:GLN:HG3	2.02	0.41
2:C:57:VAL:HB	2:C:79:VAL:HG23	2.02	0.41
2:A:122:VAL:CG2	2:A:125:ALA:HB2	2.50	0.41
2:B:157:THR:HG22	2:B:157:THR:O	2.19	0.41
2:C:207:SER:C	2:C:209:ASP:N	2.74	0.41
2:C:189:LEU:CD2	2:C:193:ARG:NH2	2.79	0.41
2:C:279:ASP:O	2:C:282:SER:HB2	2.20	0.41
2:B:63:PHE:HB2	2:B:294:SER:O	2.20	0.41
2:A:237:ASP:N	2:A:237:ASP:OD1	2.54	0.41
2:A:177:LYS:O	2:A:178:HIS:CG	2.73	0.41
2:A:264:HIS:N	2:A:264:HIS:ND1	2.67	0.41
2:B:75:SER:C	2:B:76:LEU:HD23	2.40	0.41
2:A:40:GLU:OE1	2:A:82:THR:HG21	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:106:ARG:CZ	2:B:108:GLU:OE2	2.69	0.41
2:C:149:PHE:CZ	2:C:171:PRO:HG3	2.55	0.41
2:A:167:SER:O	2:A:170:THR:CB	2.69	0.41
2:A:111:PRO:HG3	2:A:146:TRP:CD2	2.56	0.41
2:C:39:VAL:HG21	4:C:360:HOH:O	2.21	0.41
2:C:41:LEU:O	2:C:134:ARG:NH1	2.53	0.41
2:B:111:PRO:HG3	2:B:146:TRP:CE2	2.56	0.41
2:A:291:SER:OG	2:A:292:HIS:N	2.54	0.41
2:A:155:SER:O	2:A:159:GLU:HG2	2.20	0.41
2:C:110:LEU:HA	2:C:111:PRO:HD3	1.78	0.40
2:A:126:THR:HG22	2:A:137:ASN:HB2	2.01	0.40
2:A:106:ARG:HB2	2:A:108:GLU:OE1	2.21	0.40
2:B:121:ARG:HG3	2:B:121:ARG:O	2.22	0.40
2:A:303:SER:O	2:A:307:LYS:HG3	2.21	0.40
2:C:120:ILE:HG13	2:C:148:LEU:HD23	2.04	0.40
2:B:158:GLN:OE1	2:B:165:ALA:HA	2.22	0.40
2:C:64:PRO:HB2	2:C:107:PHE:CD1	2.56	0.40
2:A:72:TYR:HE2	2:A:107:PHE:CD1	2.39	0.40
2:A:232:GLN:HG3	2:A:234:HIS:CD2	2.55	0.40
2:B:238:GLU:HG3	2:B:239:TYR:CD1	2.57	0.40
2:B:84:TYR:CE2	2:B:86:LYS:HE3	2.56	0.40
1:D:6:DG:H5'	2:A:73:ILE:HD13	2.03	0.40
2:C:48:SER:C	2:C:50:GLN:N	2.73	0.40
2:B:303:SER:O	2:B:307:LYS:HG3	2.20	0.40
2:A:258:LEU:HB2	4:A:337:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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
2	A	263/320 (82%)	245 (93%)	13 (5%)	5 (2%)	10 19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	271/320 (85%)	245 (90%)	22 (8%)	4 (2%)	13	26
2	C	274/320 (86%)	242 (88%)	24 (9%)	8 (3%)	6	9
All	All	808/960 (84%)	732 (91%)	59 (7%)	17 (2%)	9	16

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	166	VAL
2	A	176	SER
2	A	178	HIS
2	C	219	ALA
2	A	186	ILE
2	C	208	SER
2	C	283	THR
2	B	49	ALA
2	C	69	GLN
2	C	144	SER
2	C	218	GLN
2	B	70	GLU
2	C	49	ALA
2	A	184	ASN
2	B	186	ILE
2	B	143	SER
2	C	205	VAL

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	
2	A	240/281 (85%)	231 (96%)	9 (4%)	40	68
2	B	246/281 (88%)	235 (96%)	11 (4%)	34	62
2	C	247/281 (88%)	235 (95%)	12 (5%)	31	57
All	All	733/843 (87%)	701 (96%)	32 (4%)	35	63

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	121	ARG
2	A	122	VAL
2	A	124	ARG
2	A	176	SER
2	A	211	TYR
2	A	240	THR
2	A	243	LEU
2	A	261	LYS
2	A	315	ASP
2	B	37	GLU
2	B	46	LEU
2	B	47	THR
2	B	121	ARG
2	B	150	SER
2	B	191	ASN
2	B	220	GLN
2	B	240	THR
2	B	243	LEU
2	B	260	LEU
2	B	266	ARG
2	C	68	ASN
2	C	80	ASP
2	C	95	SER
2	C	96	ASP
2	C	121	ARG
2	C	122	VAL
2	C	128	ARG
2	C	164	ASP
2	C	174	PHE
2	C	191	ASN
2	C	239	TYR
2	C	272	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	52	GLN
2	A	87	GLN
2	A	215	ASN
2	B	52	GLN
2	B	137	ASN

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Mol	Chain	Res	Type
2	B	162	ASN
2	B	163	GLN
2	B	178	HIS
2	B	191	ASN
2	B	234	HIS
2	B	241	ASN
2	B	251	GLN
2	C	137	ASN
2	C	241	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	A	321	-	4,4,4	0.12	0	6,6,6	0.16	0
3	SO4	A	322	-	4,4,4	0.25	0	6,6,6	0.17	0
3	SO4	B	321	-	4,4,4	0.15	0	6,6,6	0.29	0
3	SO4	B	322	-	4,4,4	0.30	0	6,6,6	0.09	0
3	SO4	C	321	-	4,4,4	0.18	0	6,6,6	0.25	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	C	322	-	4,4,4	0.30	0	6,6,6	0.23	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
3	SO4	A	321	-	-	0/0/0/0	0/0/0/0
3	SO4	A	322	-	-	0/0/0/0	0/0/0/0
3	SO4	B	321	-	-	0/0/0/0	0/0/0/0
3	SO4	B	322	-	-	0/0/0/0	0/0/0/0
3	SO4	C	321	-	-	0/0/0/0	0/0/0/0
3	SO4	C	322	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

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	D	4/6 (66%)	0.70	1 (25%) 1 0	48, 54, 60, 72	0
1	E	4/6 (66%)	0.74	1 (25%) 1 0	53, 58, 69, 75	0
2	A	269/320 (84%)	0.44	26 (9%) 10 6	22, 41, 68, 81	0
2	B	275/320 (85%)	0.52	30 (10%) 7 4	20, 44, 71, 79	0
2	C	278/320 (86%)	0.49	21 (7%) 17 12	23, 44, 69, 74	0
All	All	830/972 (85%)	0.49	79 (9%) 10 6	20, 44, 70, 81	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	86	LYS	5.8
2	C	166	VAL	5.8
2	C	94	ALA	5.6
2	B	141	PHE	5.1
2	B	142	TYR	5.0
2	A	142	TYR	4.8
2	C	178	HIS	4.5
2	B	284	GLN	4.4
2	B	49	ALA	4.3
2	B	36	TYR	4.0
2	C	180	THR	4.0
2	B	69	GLN	4.0
2	A	178	HIS	3.9
2	A	162	ASN	3.9
2	C	164	ASP	3.7
2	A	234	HIS	3.6
2	B	46	LEU	3.6
2	B	166	VAL	3.6
2	A	95	SER	3.5
2	B	168	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	163	GLN	3.4
2	C	169	THR	3.4
2	B	181	ILE	3.3
2	C	162	ASN	3.3
2	B	165	ALA	3.3
1	D	5	DG	3.3
2	A	166	VAL	3.2
2	B	178	HIS	3.2
2	C	69	GLN	3.2
2	C	163	GLN	3.1
2	B	162	ASN	3.1
2	C	234	HIS	3.1
2	A	182	GLU	2.9
2	B	180	THR	2.9
1	E	5	DG	2.9
2	A	87	GLN	2.9
2	C	280	GLU	2.8
2	B	316	ASP	2.8
2	C	95	SER	2.8
2	C	284	GLN	2.8
2	B	179	ALA	2.8
2	B	234	HIS	2.8
2	A	72	TYR	2.7
2	A	100	LEU	2.7
2	A	284	GLN	2.7
2	A	165	ALA	2.7
2	C	165	ALA	2.6
2	B	47	THR	2.6
2	C	179	ALA	2.6
2	B	105	LYS	2.6
2	B	164	ASP	2.5
2	B	176	SER	2.5
2	A	124	ARG	2.5
2	B	48	SER	2.5
2	C	316	ASP	2.5
2	C	133	GLN	2.4
2	A	130	TYR	2.3
2	C	36	TYR	2.3
2	C	88	GLN	2.3
2	C	217	ALA	2.3
2	A	107	PHE	2.3
2	A	85	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	283	THR	2.2
2	B	88	GLN	2.2
2	B	96	ASP	2.2
2	B	45	SER	2.2
2	A	181	ILE	2.2
2	A	167	SER	2.2
2	B	95	SER	2.1
2	A	39	VAL	2.1
2	A	36	TYR	2.1
2	A	37	GLU	2.1
2	A	76	LEU	2.1
2	A	43	LYS	2.1
2	B	124	ARG	2.1
2	B	58	VAL	2.0
2	A	141	PHE	2.0
2	C	218	GLN	2.0
2	A	180	THR	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	A	321	5/5	0.95	0.19	-0.04	71,73,74,75	0
3	SO4	C	322	5/5	0.98	0.17	-1.00	66,66,67,68	0
3	SO4	A	322	5/5	0.95	0.16	-1.34	81,82,83,85	0
3	SO4	B	322	5/5	0.88	0.17	-1.50	92,93,94,95	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	C	321	5/5	0.96	0.12	-1.62	71,72,73,74	0
3	SO4	B	321	5/5	0.96	0.16	-	58,59,61,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.