



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 31, 2020 – 09:07 AM BST

PDB ID : 1C0W
Title : CRYSTAL STRUCTURE OF THE COBALT-ACTIVATED DIPHTHERIA TOXIN REPRESSOR-DNA COMPLEX REVEALS A METAL BINDING SH-LIKE DOMAIN
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Deposited on : 1999-07-22
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13
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)	:	2.13

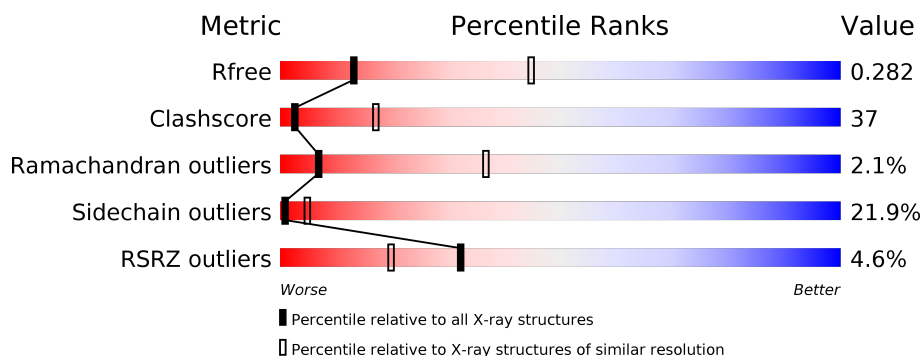
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	E	21	
2	F	21	
3	A	225	
3	B	225	
3	C	225	
3	D	225	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6575 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 DNA (5'-D(P*AP*TP*TP*AP*GP*GP*TP*TP*AP*GP*CP*CP*TP*AP*CP*CP*CP*TP*AP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	21	Total	C	N	O	P	0	0	0
			427	205	74	127	21			

- Molecule 2 is a DNA chain called DNA (5'-D(P*AP*TP*TP*AP*GP*GP*TP*TP*AP*GP*GP*CP*TP*AP*CP*CP*CP*TP*AP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	21	Total	C	N	O	P	0	0	0
			430	206	76	127	21			

- Molecule 3 is a protein called DIPHTHERIA TOXIN REPRESSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	173	Total	C	N	O	S	0	0	0
			1339	837	239	257	6			
3	B	219	Total	C	N	O	S	0	0	0
			1703	1054	305	338	6			
3	C	175	Total	C	N	O	S	0	0	0
			1335	833	241	255	6			
3	D	175	Total	C	N	O	S	0	0	0
			1333	831	241	255	6			

- Molecule 4 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Co	0	0
			2	2		
4	A	2	Total	Co	0	0
			2	2		
4	D	2	Total	Co	0	0
			2	2		

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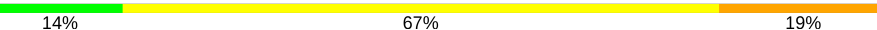
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	Co	0	0
			2	2		

3 Residue-property plots

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 (5'-D(P*AP*TP*TP*AP*GP*GP*TP*TP*AP*GP*CP*CP*TP*AP*CP*C P*CP*TP*AP*AP*T)-3')

Chain E: 

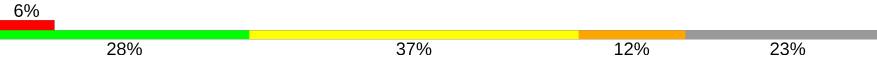


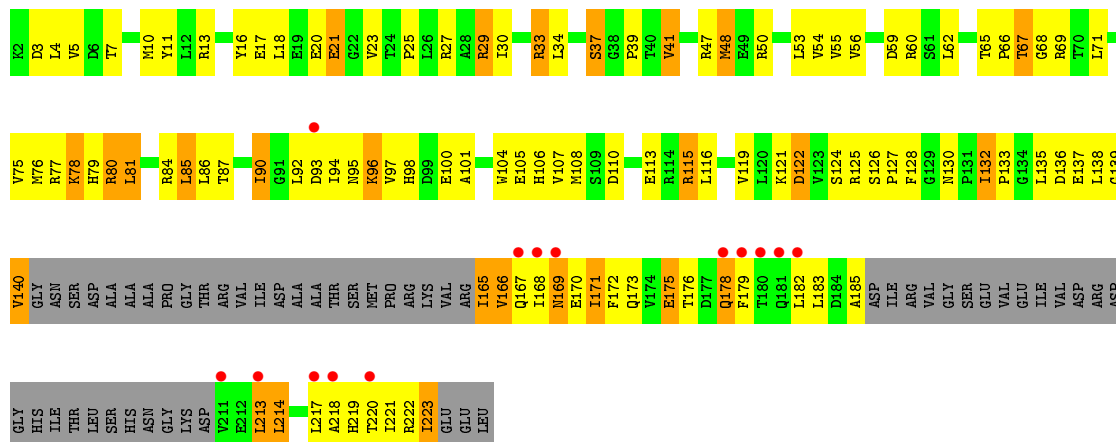
- Molecule 2: DNA (5'-D(P*AP*TP*TP*AP*GP*GP*TP*TP*AP*GP*GP*CP*TP*AP*CP*C P*CP*TP*AP*AP*T)-3')

Chain F: 

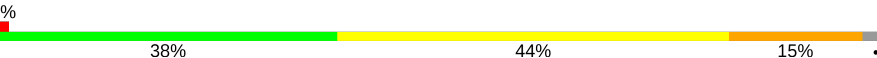


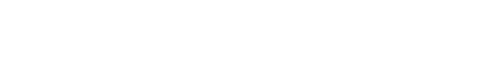
- Molecule 3: DIPHTHERIA TOXIN REPRESSOR

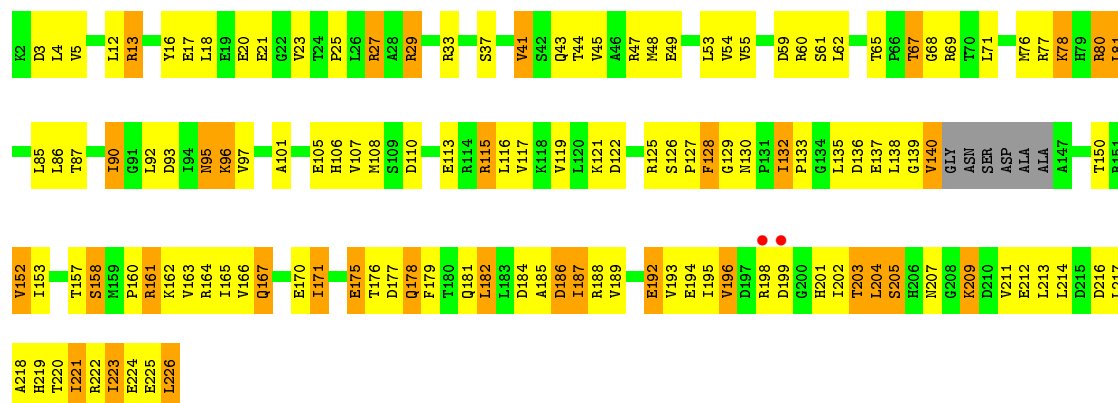
Chain A: 



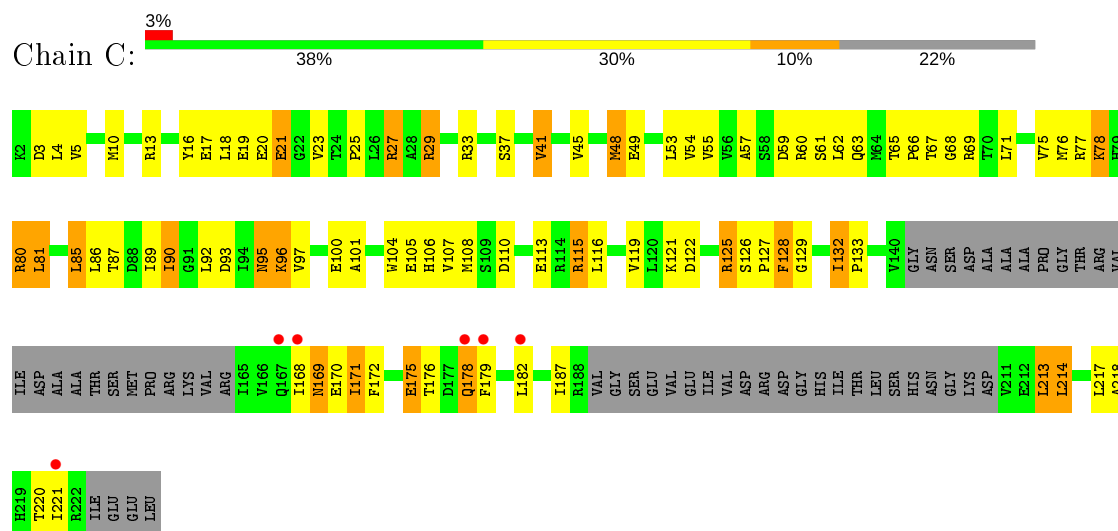
- Molecule 3: DIPHTHERIA TOXIN REPRESSOR

Chain B: 

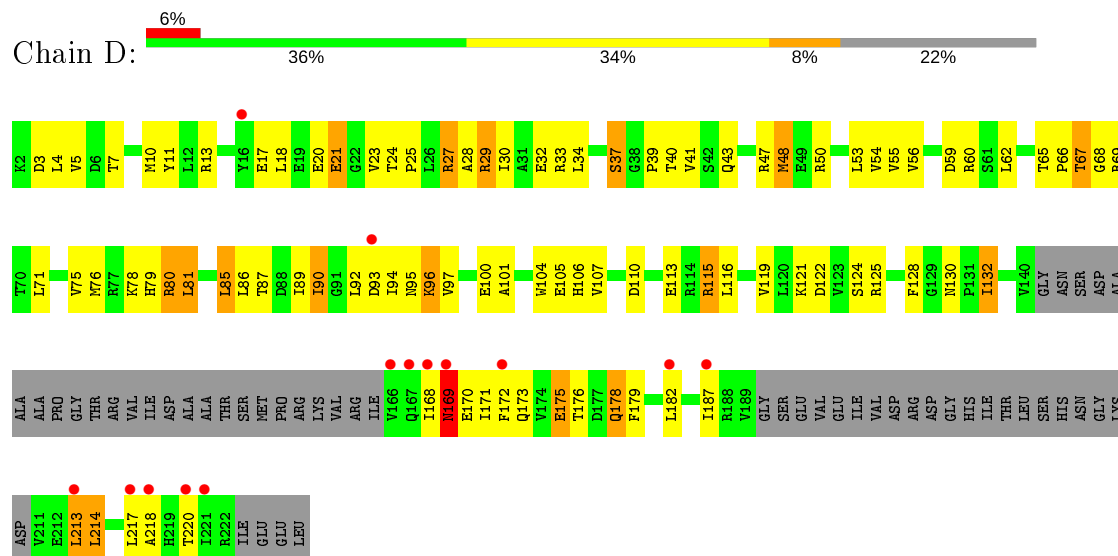




- Molecule 3: DIPHTHERIA TOXIN REPRESSOR



- Molecule 3: DIPHTHERIA TOXIN REPRESSOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 ₂ 2 ₂	Depositor
Cell constants a, b, c, α , β , γ	180.20 Å 180.20 Å 151.00 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.20 48.69 – 3.18	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-3.20) 99.2 (48.69-3.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.95 (at 3.19 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.249 , 0.279 0.254 , 0.282	Depositor DCC
R_{free} test set	2105 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 100.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	6575	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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:
CO

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	E	0.67	0/477	0.87	0/733
2	F	0.64	0/481	0.84	0/740
3	A	0.45	0/1353	0.65	0/1835
3	B	0.43	0/1722	0.68	0/2335
3	C	0.39	0/1349	0.60	0/1830
3	D	0.39	0/1347	0.62	0/1827
All	All	0.46	0/6729	0.68	0/9300

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	E	0	4
2	F	0	2
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	405	DG	Sidechain
1	E	406	DG	Sidechain
1	E	411	DC	Sidechain
1	E	416	DC	Sidechain
2	F	511	DG	Sidechain

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Mol	Chain	Res	Type	Group
2	F	516	DC	Sidechain

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	E	427	0	239	36	0
2	F	430	0	239	35	0
3	A	1339	0	1338	101	0
3	B	1703	0	1698	141	0
3	C	1335	0	1313	84	0
3	D	1333	0	1305	89	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
All	All	6575	0	6132	464	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 37.

All (464) 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)
3:A:165:ILE:HA	3:A:223:ILE:HG13	1.29	1.09
1:E:420:DA:H2''	1:E:421:DT:H5''	1.38	1.04
2:F:506:DG:H2''	2:F:507:DT:H5'	1.34	1.02
2:F:506:DG:H2''	2:F:507:DT:C5'	1.89	1.02
3:D:175:GLU:HG3	3:D:178:GLN:HB3	1.39	1.02
3:C:175:GLU:HG3	3:C:178:GLN:HB3	1.42	1.01
3:A:166:VAL:HB	3:A:223:ILE:HA	1.42	1.00
3:A:175:GLU:HG3	3:A:178:GLN:HB3	1.45	0.97
3:A:165:ILE:CA	3:A:223:ILE:HG13	1.95	0.96
3:B:161:ARG:HB2	3:B:161:ARG:HH11	1.30	0.93
3:A:172:PHE:HB2	3:A:220:THR:HG21	1.52	0.92
3:C:172:PHE:HB2	3:C:220:THR:CG2	2.00	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:172:PHE:HB2	3:D:220:THR:HG21	1.52	0.92
1:E:417:DC:H2'	1:E:418:DT:H71	1.56	0.88
3:B:214:LEU:HB2	3:B:217:LEU:HD23	1.57	0.85
3:C:172:PHE:HB2	3:C:220:THR:HG21	1.56	0.85
3:C:175:GLU:HG2	3:C:217:LEU:HD11	1.57	0.84
3:D:172:PHE:HB2	3:D:220:THR:CG2	2.08	0.84
3:B:87:THR:HG22	3:B:97:VAL:HG21	1.61	0.83
3:A:172:PHE:HB2	3:A:220:THR:CG2	2.09	0.83
1:E:414:DA:H2''	1:E:415:DC:C5'	2.09	0.82
3:D:175:GLU:CG	3:D:178:GLN:HB3	2.09	0.82
3:D:101:ALA:O	3:D:105:GLU:HG2	1.80	0.82
1:E:401:DA:OP2	3:D:28:ALA:HB2	1.81	0.81
1:E:403:DT:H2''	1:E:404:DA:H5'	1.64	0.80
3:A:175:GLU:CG	3:A:178:GLN:HB3	2.11	0.79
3:A:94:ILE:HD11	3:A:128:PHE:HE1	1.46	0.79
1:E:414:DA:H2''	1:E:415:DC:H5''	1.64	0.79
2:F:507:DT:H5'	2:F:507:DT:H6	1.48	0.79
3:D:94:ILE:HD11	3:D:128:PHE:HE1	1.48	0.78
3:B:165:ILE:HG22	3:B:189:VAL:HA	1.64	0.78
3:A:165:ILE:HD11	3:A:221:ILE:HG22	1.67	0.77
3:C:169:ASN:HB2	3:C:220:THR:O	1.84	0.76
3:B:165:ILE:CG2	3:B:189:VAL:HA	2.16	0.75
3:A:115:ARG:O	3:A:119:VAL:HG23	1.86	0.75
3:D:213:LEU:HD12	3:D:218:ALA:HA	1.68	0.74
1:E:420:DA:C2'	1:E:421:DT:H5''	2.16	0.74
3:B:184:ASP:C	3:B:186:ASP:H	1.91	0.74
3:C:115:ARG:HD3	3:D:92:LEU:HA	1.69	0.74
1:E:406:DG:H5''	3:A:60:ARG:HD3	1.70	0.74
3:B:164:ARG:NH1	3:B:226:LEU:HA	2.02	0.73
1:E:401:DA:H2'	1:E:402:DT:C7	2.19	0.73
2:F:506:DG:H2'	2:F:507:DT:H72	1.69	0.73
2:F:507:DT:H2'	2:F:508:DT:H71	1.71	0.73
1:E:414:DA:C2'	1:E:415:DC:H5''	2.19	0.72
3:C:92:LEU:HA	3:D:115:ARG:HD3	1.70	0.72
2:F:505:DG:H1'	2:F:506:DG:H5''	1.72	0.72
3:A:65:THR:OG1	3:A:67:THR:HG23	1.91	0.71
3:A:166:VAL:HG13	3:A:167:GLN:HG2	1.71	0.71
3:B:128:PHE:HD1	3:B:128:PHE:N	1.89	0.70
3:C:81:LEU:HD13	3:C:132:ILE:HG21	1.74	0.70
2:F:512:DC:H2'	2:F:513:DT:H71	1.74	0.70
2:F:516:DC:H2''	2:F:517:DC:OP2	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:161:ARG:NH1	3:B:161:ARG:HB2	2.05	0.70
1:E:417:DC:C2'	1:E:418:DT:H71	2.22	0.70
3:C:171:ILE:HG22	3:C:220:THR:HG23	1.74	0.70
2:F:505:DG:C2'	2:F:506:DG:H5''	2.22	0.70
3:A:101:ALA:O	3:A:105:GLU:HG2	1.90	0.69
3:C:172:PHE:HB2	3:C:220:THR:HG22	1.73	0.69
3:B:128:PHE:CD1	3:B:128:PHE:N	2.60	0.69
3:A:87:THR:HG22	3:A:97:VAL:HG21	1.74	0.69
3:C:93:ASP:HB3	3:C:96:LYS:HG3	1.75	0.69
3:A:175:GLU:HG2	3:A:217:LEU:HD11	1.75	0.68
3:B:93:ASP:HB3	3:B:96:LYS:HG3	1.74	0.68
3:D:115:ARG:O	3:D:119:VAL:HG23	1.94	0.68
3:C:101:ALA:O	3:C:105:GLU:HG2	1.93	0.68
1:E:406:DG:H4'	3:A:60:ARG:CZ	2.23	0.68
3:B:203:THR:HG23	3:B:212:GLU:HA	1.76	0.68
3:A:93:ASP:HB3	3:A:96:LYS:HG3	1.75	0.67
3:C:87:THR:HG22	3:C:97:VAL:HG21	1.77	0.67
3:B:163:VAL:HG23	3:B:224:GLU:O	1.95	0.67
3:D:93:ASP:HB3	3:D:96:LYS:HG3	1.77	0.67
1:E:405:DG:H1'	1:E:406:DG:H5'	1.77	0.67
3:B:193:VAL:HG11	3:B:204:LEU:HB3	1.77	0.66
3:B:224:GLU:HG2	3:B:225:GLU:H	1.58	0.66
3:D:87:THR:HG22	3:D:97:VAL:HG21	1.77	0.66
3:B:166:VAL:HG13	3:B:224:GLU:HB2	1.77	0.66
3:D:71:LEU:O	3:D:75:VAL:HG23	1.96	0.66
1:E:401:DA:H2'	1:E:402:DT:H72	1.77	0.66
3:A:115:ARG:HD3	3:B:92:LEU:HA	1.77	0.66
2:F:517:DC:H2'	2:F:518:DT:H72	1.78	0.66
3:B:115:ARG:O	3:B:119:VAL:HG23	1.96	0.65
3:B:166:VAL:CG1	3:B:224:GLU:HB2	2.26	0.65
3:C:86:LEU:HD22	3:C:92:LEU:HD22	1.76	0.65
3:C:78:LYS:HE3	3:C:108:MET:O	1.97	0.65
1:E:414:DA:H2''	1:E:415:DC:H5'	1.79	0.65
3:C:27:ARG:HD3	3:C:62:LEU:HD21	1.78	0.65
3:D:169:ASN:HB2	3:D:220:THR:O	1.97	0.65
3:B:162:LYS:O	3:B:226:LEU:HD11	1.96	0.65
3:B:165:ILE:HA	3:B:223:ILE:CD1	2.26	0.64
3:B:166:VAL:HG22	3:B:223:ILE:HA	1.78	0.64
3:C:115:ARG:O	3:C:119:VAL:HG23	1.98	0.64
3:C:80:ARG:HB2	3:C:132:ILE:HD13	1.79	0.64
3:D:100:GLU:HG3	3:D:104:TRP:NE1	2.13	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:163:VAL:CG2	3:B:223:ILE:HG13	2.28	0.63
2:F:505:DG:H2''	2:F:506:DG:H5''	1.78	0.63
3:C:128:PHE:N	3:C:128:PHE:HD1	1.97	0.63
3:A:10:MET:SD	3:A:106:HIS:CE1	2.92	0.63
3:D:86:LEU:HD22	3:D:92:LEU:HD22	1.81	0.62
3:D:10:MET:SD	3:D:106:HIS:CE1	2.92	0.62
3:B:163:VAL:HG22	3:B:223:ILE:HG13	1.81	0.62
3:D:11:TYR:HB3	3:D:48:MET:HE1	1.81	0.61
3:B:162:LYS:NZ	3:B:192:GLU:HG2	2.16	0.61
3:A:166:VAL:CB	3:A:223:ILE:HA	2.26	0.61
3:B:164:ARG:HH12	3:B:226:LEU:HA	1.65	0.61
3:C:172:PHE:CB	3:C:220:THR:HG21	2.29	0.61
3:C:168:ILE:HG22	3:C:172:PHE:HB3	1.83	0.61
3:C:175:GLU:CG	3:C:178:GLN:HB3	2.26	0.61
2:F:506:DG:H2''	2:F:507:DT:H5''	1.80	0.61
3:B:184:ASP:C	3:B:186:ASP:N	2.53	0.61
3:B:207:ASN:C	3:B:209:LYS:H	2.05	0.60
3:B:4:LEU:O	3:B:5:VAL:HB	1.99	0.60
3:A:165:ILE:HD12	3:A:223:ILE:HB	1.83	0.60
3:B:164:ARG:CZ	3:B:226:LEU:HA	2.31	0.60
3:A:166:VAL:HG12	3:A:222:ARG:C	2.22	0.60
1:E:411:DC:H2''	1:E:412:DC:O5'	2.02	0.60
3:C:213:LEU:HD12	3:C:218:ALA:HA	1.82	0.59
3:C:214:LEU:HG	3:C:217:LEU:HD12	1.84	0.59
3:D:4:LEU:HD12	3:D:7:THR:HG23	1.84	0.59
3:A:86:LEU:HA	3:A:90:ILE:HG13	1.84	0.59
1:E:407:DT:H2'	1:E:408:DT:H72	1.84	0.59
3:C:85:LEU:HD22	3:C:116:LEU:HD21	1.83	0.59
3:A:54:VAL:HG22	3:A:55:VAL:N	2.17	0.59
3:B:138:LEU:HD23	3:B:140:VAL:HG21	1.85	0.59
3:D:65:THR:HB	3:D:66:PRO:HD2	1.83	0.59
3:B:27:ARG:HD3	3:B:62:LEU:HD21	1.83	0.59
3:C:128:PHE:N	3:C:128:PHE:CD1	2.68	0.59
3:D:128:PHE:HD2	3:D:173:GLN:CD	2.06	0.59
3:A:128:PHE:HD2	3:A:173:GLN:CD	2.05	0.58
3:A:71:LEU:O	3:A:75:VAL:HG23	2.03	0.58
3:A:94:ILE:HD11	3:A:128:PHE:CE1	2.34	0.58
3:D:29:ARG:NH1	3:D:32:GLU:CD	2.56	0.58
1:E:401:DA:P	3:D:27:ARG:HB2	2.43	0.58
1:E:415:DC:H2''	1:E:416:DC:H5'	1.86	0.58
2:F:507:DT:H5'	2:F:507:DT:C6	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:168:ILE:HG22	3:D:172:PHE:HB3	1.85	0.57
3:A:135:LEU:O	3:A:140:VAL:HG23	2.03	0.57
3:A:213:LEU:HD12	3:A:218:ALA:HA	1.85	0.57
3:D:81:LEU:HD13	3:D:132:ILE:HG21	1.86	0.57
3:B:54:VAL:HG22	3:B:55:VAL:N	2.18	0.57
3:B:5:VAL:HG11	3:B:106:HIS:HB2	1.86	0.57
3:B:153:ILE:HA	3:B:202:ILE:CD1	2.34	0.57
3:B:203:THR:CG2	3:B:212:GLU:HA	2.35	0.57
3:A:33:ARG:CZ	3:A:171:ILE:HG13	2.35	0.57
3:A:4:LEU:O	3:A:5:VAL:HB	2.03	0.57
3:C:178:GLN:HG2	3:C:217:LEU:HD13	1.87	0.57
3:D:113:GLU:O	3:D:116:LEU:HB2	2.05	0.57
3:B:65:THR:OG1	3:B:67:THR:HG23	2.04	0.56
3:C:54:VAL:HG22	3:C:55:VAL:N	2.20	0.56
1:E:417:DC:H2'	1:E:418:DT:C7	2.33	0.56
3:B:171:ILE:HD13	3:B:171:ILE:O	2.05	0.56
2:F:505:DG:C1'	2:F:506:DG:H5''	2.35	0.56
3:A:29:ARG:HH11	3:A:29:ARG:HA	1.70	0.56
3:B:164:ARG:O	3:B:223:ILE:HD12	2.06	0.56
3:C:10:MET:SD	3:C:106:HIS:CE1	2.98	0.56
3:B:205:SER:HA	3:B:209:LYS:O	2.05	0.56
2:F:506:DG:H2'	2:F:507:DT:C7	2.36	0.56
1:E:416:DC:H2''	1:E:417:DC:O5'	2.05	0.56
2:F:501:DA:OP1	3:B:60:ARG:HB3	2.05	0.56
3:A:92:LEU:HA	3:B:115:ARG:HD3	1.87	0.56
3:D:86:LEU:HA	3:D:90:ILE:HG13	1.88	0.55
3:B:41:VAL:O	3:B:45:VAL:HG23	2.06	0.55
3:D:11:TYR:HB3	3:D:48:MET:CE	2.36	0.55
3:B:80:ARG:HB2	3:B:132:ILE:HD13	1.88	0.55
3:B:158:SER:O	3:B:198:ARG:NH2	2.40	0.55
3:B:29:ARG:HA	3:B:29:ARG:HH11	1.71	0.55
2:F:505:DG:H2''	2:F:506:DG:C5'	2.37	0.55
3:C:93:ASP:HB3	3:C:96:LYS:CG	2.36	0.55
3:D:30:ILE:HG23	3:D:34:LEU:HD12	1.87	0.55
3:C:71:LEU:O	3:C:75:VAL:HG23	2.07	0.55
3:B:160:PRO:CA	3:B:196:VAL:HG13	2.36	0.54
3:C:182:LEU:O	3:C:187:ILE:HG12	2.06	0.54
3:A:81:LEU:HD13	3:A:132:ILE:HG21	1.88	0.54
1:E:403:DT:C2'	1:E:404:DA:H5'	2.35	0.54
3:D:93:ASP:O	3:D:97:VAL:HG23	2.08	0.54
3:B:59:ASP:OD1	3:B:61:SER:HB3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:178:GLN:HG3	3:C:178:GLN:O	2.06	0.54
3:D:214:LEU:HG	3:D:217:LEU:HD12	1.89	0.54
3:B:166:VAL:HG22	3:B:223:ILE:CA	2.37	0.54
3:A:214:LEU:HG	3:A:217:LEU:HD12	1.88	0.53
3:C:80:ARG:HH21	3:C:170:GLU:HG2	1.73	0.53
3:D:48:MET:HB3	3:D:54:VAL:HG12	1.90	0.53
3:B:195:ILE:CG2	3:B:204:LEU:HD23	2.39	0.53
3:A:113:GLU:O	3:A:116:LEU:HB2	2.08	0.53
3:B:162:LYS:HZ2	3:B:192:GLU:HG2	1.73	0.53
3:D:76:MET:HE1	3:D:79:HIS:CD2	2.43	0.53
3:A:105:GLU:HG3	3:A:106:HIS:CD2	2.44	0.53
3:B:193:VAL:HG12	3:B:194:GLU:N	2.23	0.53
3:A:168:ILE:HG22	3:A:172:PHE:HB3	1.89	0.53
3:D:17:GLU:O	3:D:21:GLU:HG2	2.09	0.52
3:A:16:TYR:CE1	3:A:76:MET:HG2	2.44	0.52
1:E:401:DA:OP2	3:D:28:ALA:CB	2.56	0.52
2:F:517:DC:C6	2:F:518:DT:H72	2.45	0.52
3:C:93:ASP:O	3:C:97:VAL:HG23	2.10	0.52
3:D:54:VAL:HG22	3:D:55:VAL:N	2.25	0.52
3:A:178:GLN:HG3	3:A:178:GLN:O	2.09	0.52
3:C:80:ARG:HH21	3:C:170:GLU:CG	2.22	0.52
3:B:203:THR:HG22	3:B:211:VAL:O	2.10	0.52
3:A:53:LEU:O	3:A:68:GLY:HA3	2.10	0.52
3:B:87:THR:CG2	3:B:97:VAL:HG21	2.38	0.52
3:C:4:LEU:O	3:C:5:VAL:HB	2.09	0.52
3:D:89:ILE:N	3:D:89:ILE:HD12	2.24	0.52
2:F:517:DC:H2'	2:F:518:DT:C7	2.40	0.52
3:B:81:LEU:HD13	3:B:132:ILE:HG21	1.92	0.52
3:D:213:LEU:CD1	3:D:218:ALA:HA	2.39	0.52
3:A:87:THR:CG2	3:A:97:VAL:HG21	2.40	0.51
3:B:12:LEU:CD1	3:B:71:LEU:HD23	2.41	0.51
3:B:135:LEU:O	3:B:140:VAL:HG23	2.10	0.51
3:B:53:LEU:O	3:B:68:GLY:HA3	2.11	0.51
3:D:29:ARG:HH11	3:D:32:GLU:CD	2.13	0.51
3:B:93:ASP:HB3	3:B:96:LYS:CG	2.38	0.51
3:C:59:ASP:O	3:C:60:ARG:HB2	2.10	0.51
3:A:183:LEU:C	3:A:185:ALA:H	2.13	0.51
3:B:164:ARG:NH2	3:B:226:LEU:HA	2.25	0.51
3:B:185:ALA:HB3	3:B:187:ILE:HG13	1.93	0.50
3:A:165:ILE:HD11	3:A:221:ILE:CG2	2.38	0.50
3:A:98:HIS:CE1	3:A:170:GLU:OE1	2.64	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:128:PHE:N	3:D:128:PHE:CD1	2.79	0.50
2:F:516:DC:H2'	3:D:40:THR:CG2	2.41	0.50
3:B:182:LEU:HG	3:B:187:ILE:HB	1.94	0.50
3:B:165:ILE:HA	3:B:223:ILE:HD11	1.93	0.50
3:A:100:GLU:HG3	3:A:104:TRP:NE1	2.27	0.50
3:A:128:PHE:CD1	3:A:128:PHE:N	2.79	0.50
3:B:160:PRO:HB3	3:B:196:VAL:HG13	1.93	0.50
3:B:165:ILE:HD12	3:B:223:ILE:HD11	1.94	0.50
3:D:182:LEU:O	3:D:187:ILE:HG12	2.11	0.50
3:B:49:GLU:HA	3:B:54:VAL:O	2.12	0.50
3:D:178:GLN:HG3	3:D:178:GLN:O	2.12	0.50
3:B:13:ARG:O	3:B:16:TYR:HB3	2.12	0.50
1:E:417:DC:H2''	1:E:418:DT:C6	2.46	0.50
3:C:168:ILE:CG2	3:C:172:PHE:HB3	2.42	0.49
3:B:153:ILE:HG22	3:B:219:HIS:HA	1.94	0.49
3:C:16:TYR:CE1	3:C:76:MET:HG2	2.46	0.49
3:B:223:ILE:HG23	3:B:224:GLU:N	2.28	0.49
2:F:513:DT:H2''	2:F:514:DA:C8	2.47	0.49
3:C:179:PHE:O	3:C:182:LEU:HB3	2.11	0.49
3:A:21:GLU:HB2	3:A:23:VAL:HG23	1.94	0.49
3:D:168:ILE:CG2	3:D:172:PHE:HB3	2.43	0.49
3:A:178:GLN:OE1	3:A:213:LEU:HA	2.13	0.49
3:A:48:MET:HB3	3:A:54:VAL:HG12	1.95	0.49
3:C:41:VAL:O	3:C:45:VAL:HG23	2.13	0.49
3:D:80:ARG:CZ	3:D:130:ASN:HB3	2.42	0.49
3:B:160:PRO:CB	3:B:196:VAL:HG13	2.43	0.48
3:B:163:VAL:HG21	3:B:223:ILE:CG2	2.43	0.48
3:B:150:THR:O	3:B:223:ILE:N	2.47	0.48
3:D:37:SER:OG	3:D:39:PRO:HD2	2.14	0.48
1:E:406:DG:H5''	3:A:60:ARG:CD	2.43	0.48
3:A:37:SER:OG	3:A:39:PRO:HD2	2.13	0.48
3:B:195:ILE:HG22	3:B:204:LEU:HD23	1.95	0.48
3:A:5:VAL:HG11	3:A:106:HIS:HB2	1.96	0.48
3:A:65:THR:HB	3:A:66:PRO:HD2	1.94	0.48
3:A:93:ASP:HB3	3:A:96:LYS:CG	2.44	0.48
3:B:80:ARG:HD3	3:B:126:SER:HB3	1.96	0.48
3:A:169:ASN:HB2	3:A:220:THR:O	2.13	0.48
3:C:29:ARG:HH11	3:C:29:ARG:HA	1.79	0.48
3:B:113:GLU:O	3:B:117:VAL:HG23	2.14	0.48
3:B:76:MET:HE2	3:B:170:GLU:HG3	1.95	0.48
3:A:166:VAL:HG13	3:A:167:GLN:CG	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:5:VAL:HG11	3:C:106:HIS:HB2	1.96	0.48
3:C:48:MET:HB3	3:C:54:VAL:HG12	1.96	0.48
3:B:80:ARG:CZ	3:B:130:ASN:HB3	2.43	0.47
3:D:94:ILE:HD11	3:D:128:PHE:CE1	2.38	0.47
3:B:217:LEU:N	3:B:217:LEU:HD22	2.28	0.47
3:D:18:LEU:HD13	3:D:25:PRO:HA	1.95	0.47
3:D:47:ARG:HG3	3:D:50:ARG:NH2	2.29	0.47
3:B:21:GLU:OE2	3:B:171:ILE:HB	2.15	0.47
3:C:76:MET:O	3:C:80:ARG:HG3	2.15	0.47
3:B:160:PRO:C	3:B:161:ARG:HG3	2.34	0.47
3:C:172:PHE:HE1	3:C:179:PHE:HA	1.79	0.47
3:A:110:ASP:O	3:A:113:GLU:HB2	2.14	0.47
3:A:76:MET:O	3:A:80:ARG:HG3	2.15	0.47
3:B:175:GLU:HB3	3:B:178:GLN:HB3	1.95	0.47
3:B:193:VAL:CG1	3:B:194:GLU:N	2.77	0.47
3:C:17:GLU:O	3:C:20:GLU:HB2	2.14	0.47
1:E:414:DA:H1'	1:E:415:DC:H5''	1.96	0.47
3:B:113:GLU:O	3:B:116:LEU:HB2	2.14	0.47
1:E:412:DC:H2'	1:E:413:DT:H72	1.96	0.47
2:F:517:DC:C2'	2:F:518:DT:H72	2.44	0.47
3:A:86:LEU:HD22	3:A:92:LEU:HD22	1.95	0.47
3:B:223:ILE:HG23	3:B:224:GLU:O	2.15	0.47
3:C:54:VAL:N	3:C:65:THR:HG23	2.30	0.46
3:C:86:LEU:HA	3:C:90:ILE:HG13	1.96	0.46
3:B:21:GLU:HB2	3:B:23:VAL:HG23	1.97	0.46
2:F:506:DG:C2'	2:F:507:DT:C5'	2.79	0.46
3:A:80:ARG:HB3	3:A:126:SER:HB3	1.97	0.46
2:F:512:DC:C2'	2:F:513:DT:H71	2.44	0.46
3:A:183:LEU:C	3:A:185:ALA:N	2.68	0.46
3:C:81:LEU:CD1	3:C:132:ILE:HG21	2.43	0.46
3:D:85:LEU:HD12	3:D:85:LEU:O	2.16	0.46
3:B:175:GLU:O	3:B:179:PHE:N	2.43	0.46
3:B:203:THR:HG23	3:B:212:GLU:HG3	1.97	0.46
3:C:182:LEU:HD22	3:C:187:ILE:HG13	1.98	0.46
2:F:502:DT:C6	2:F:503:DT:H72	2.51	0.46
3:C:127:PRO:C	3:C:128:PHE:HD1	2.18	0.46
2:F:516:DC:H5	3:D:43:GLN:HE22	1.62	0.46
3:C:171:ILE:HG22	3:C:220:THR:CG2	2.43	0.46
3:C:53:LEU:C	3:C:65:THR:HG23	2.36	0.46
3:C:21:GLU:HB2	3:C:23:VAL:HG23	1.98	0.46
1:E:413:DT:H2''	1:E:414:DA:C8	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:11:TYR:OH	3:A:41:VAL:HA	2.16	0.46
3:A:54:VAL:CG2	3:A:55:VAL:N	2.79	0.46
3:B:158:SER:O	3:B:196:VAL:HG12	2.16	0.46
3:C:107:VAL:HG21	3:D:107:VAL:HG21	1.97	0.46
3:B:164:ARG:HH22	3:B:226:LEU:C	2.19	0.45
3:C:87:THR:CG2	3:C:97:VAL:HG21	2.45	0.45
3:B:44:THR:HA	3:B:47:ARG:NH2	2.31	0.45
3:B:54:VAL:N	3:B:65:THR:HG23	2.31	0.45
3:C:21:GLU:OE2	3:C:171:ILE:HB	2.16	0.45
3:D:28:ALA:O	3:D:32:GLU:HG2	2.16	0.45
2:F:511:DG:H2"	2:F:512:DC:OP2	2.16	0.45
2:F:501:DA:O5'	3:B:27:ARG:NH2	2.49	0.45
3:C:172:PHE:CE2	3:C:221:ILE:HD11	2.52	0.45
3:D:130:ASN:OD1	3:D:173:GLN:NE2	2.49	0.45
3:A:54:VAL:N	3:A:65:THR:HG23	2.32	0.45
3:B:178:GLN:OE1	3:B:213:LEU:HA	2.17	0.45
3:B:195:ILE:HD12	3:B:195:ILE:O	2.17	0.45
3:A:76:MET:HE1	3:A:79:HIS:CD2	2.52	0.45
3:B:78:LYS:HE3	3:B:108:MET:O	2.16	0.45
3:A:122:ASP:OD2	3:A:124:SER:HB2	2.17	0.45
3:B:136:ASP:O	3:B:139:GLY:N	2.50	0.45
3:B:225:GLU:OE1	3:B:225:GLU:HA	2.16	0.45
3:B:59:ASP:O	3:B:60:ARG:HB2	2.16	0.45
3:C:182:LEU:HD13	3:C:187:ILE:HB	1.99	0.45
3:D:17:GLU:O	3:D:20:GLU:HB2	2.16	0.45
1:E:401:DA:C2	1:E:402:DT:C2	3.05	0.45
3:C:80:ARG:HB3	3:C:126:SER:HB3	1.99	0.45
3:D:179:PHE:O	3:D:182:LEU:HB3	2.17	0.45
3:D:4:LEU:O	3:D:5:VAL:HB	2.17	0.45
3:A:179:PHE:O	3:A:182:LEU:HB3	2.16	0.45
3:B:207:ASN:C	3:B:209:LYS:N	2.69	0.45
3:D:23:VAL:HG12	3:D:24:THR:N	2.31	0.45
3:A:130:ASN:OD1	3:A:173:GLN:NE2	2.49	0.45
3:B:160:PRO:HB3	3:B:196:VAL:HG22	1.99	0.45
3:B:184:ASP:O	3:B:186:ASP:N	2.50	0.45
3:A:56:VAL:HG22	3:A:62:LEU:CD2	2.46	0.44
3:D:54:VAL:N	3:D:65:THR:HG23	2.33	0.44
3:A:84:ARG:HD2	3:A:127:PRO:HG3	1.99	0.44
3:B:166:VAL:HG23	3:B:222:ARG:C	2.37	0.44
3:B:29:ARG:HA	3:B:29:ARG:HD3	1.79	0.44
3:A:213:LEU:CD1	3:A:218:ALA:HA	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:101:ALA:O	3:B:105:GLU:HG2	2.16	0.44
3:A:169:ASN:O	3:A:173:GLN:HG3	2.18	0.44
3:B:93:ASP:O	3:B:97:VAL:HG23	2.17	0.44
3:C:53:LEU:O	3:C:68:GLY:HA3	2.17	0.44
3:B:171:ILE:CG2	3:B:220:THR:HG21	2.48	0.44
3:B:203:THR:HG22	3:B:211:VAL:C	2.38	0.44
3:D:217:LEU:O	3:D:220:THR:HB	2.17	0.44
3:B:165:ILE:HG21	3:B:189:VAL:HA	1.99	0.44
3:B:193:VAL:CG1	3:B:204:LEU:HB3	2.46	0.44
3:B:163:VAL:HG21	3:B:223:ILE:HG13	1.99	0.44
2:F:515:DC:C2	2:F:516:DC:C5	3.06	0.44
2:F:516:DC:H2'	3:D:40:THR:HG21	1.99	0.44
3:A:166:VAL:HG12	3:A:222:ARG:O	2.17	0.43
3:A:18:LEU:HD13	3:A:25:PRO:HA	2.00	0.43
3:B:165:ILE:O	3:B:165:ILE:HG22	2.18	0.43
3:C:100:GLU:HG3	3:C:104:TRP:NE1	2.33	0.43
3:D:53:LEU:O	3:D:68:GLY:HA3	2.18	0.43
3:A:4:LEU:HD12	3:A:7:THR:HG23	2.00	0.43
3:C:16:TYR:O	3:C:19:GLU:HB2	2.19	0.43
3:D:56:VAL:HG22	3:D:62:LEU:CD2	2.48	0.43
2:F:517:DC:H2''	2:F:518:DT:C6	2.54	0.43
3:A:138:LEU:HA	3:A:138:LEU:HD12	1.71	0.43
3:A:33:ARG:NH1	3:A:171:ILE:HG13	2.33	0.43
3:C:104:TRP:CD1	3:C:104:TRP:N	2.86	0.43
1:E:416:DC:H5	3:B:43:GLN:HE22	1.67	0.43
3:B:127:PRO:HB2	3:B:128:PHE:CD1	2.54	0.43
3:C:81:LEU:HD12	3:C:81:LEU:HA	1.83	0.43
3:D:172:PHE:HE1	3:D:179:PHE:HA	1.84	0.43
3:A:30:ILE:HG23	3:A:34:LEU:HD12	2.01	0.43
1:E:401:DA:P	3:D:27:ARG:CB	3.07	0.43
3:B:166:VAL:HG23	3:B:222:ARG:O	2.19	0.43
3:C:59:ASP:OD1	3:C:61:SER:HB3	2.18	0.43
3:C:95:ASN:N	3:C:95:ASN:OD1	2.52	0.43
3:D:81:LEU:HA	3:D:81:LEU:HD12	1.91	0.43
3:A:165:ILE:HG23	3:A:165:ILE:O	2.17	0.43
3:A:172:PHE:HB2	3:A:220:THR:HG22	1.95	0.43
3:A:168:ILE:CG2	3:A:172:PHE:HB3	2.49	0.43
3:A:77:ARG:HA	3:A:133:PRO:HD2	2.01	0.43
3:B:81:LEU:HD12	3:B:81:LEU:HA	1.88	0.43
3:B:86:LEU:HD22	3:B:92:LEU:HD22	2.00	0.43
3:D:4:LEU:CD1	3:D:7:THR:HG23	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:96:LYS:H	3:A:96:LYS:HG2	1.51	0.43
3:D:86:LEU:HD22	3:D:92:LEU:CD2	2.48	0.43
3:A:104:TRP:N	3:A:104:TRP:CD1	2.87	0.42
3:B:163:VAL:HG21	3:B:223:ILE:HG23	2.01	0.42
3:C:214:LEU:HD23	3:C:214:LEU:N	2.33	0.42
3:D:96:LYS:HG2	3:D:96:LYS:H	1.50	0.42
3:A:84:ARG:HD2	3:A:127:PRO:CD	2.49	0.42
3:C:125:ARG:HD3	3:C:129:GLY:HA2	2.01	0.42
3:D:65:THR:OG1	3:D:67:THR:HG23	2.19	0.42
1:E:401:DA:H2''	1:E:402:DT:O5'	2.19	0.42
3:A:47:ARG:HG3	3:A:50:ARG:NH2	2.34	0.42
3:B:198:ARG:NE	3:B:198:ARG:HA	2.34	0.42
3:B:86:LEU:HA	3:B:90:ILE:HG13	2.00	0.42
3:D:128:PHE:CD2	3:D:173:GLN:CD	2.90	0.42
3:D:175:GLU:HG2	3:D:217:LEU:HD11	2.01	0.42
2:F:507:DT:H2'	2:F:508:DT:C7	2.46	0.42
3:D:218:ALA:C	3:D:220:THR:H	2.22	0.42
3:D:27:ARG:HD3	3:D:62:LEU:HD21	2.01	0.42
3:A:115:ARG:HB3	3:A:115:ARG:HE	1.59	0.42
3:C:18:LEU:HD13	3:C:25:PRO:HA	2.02	0.42
3:D:10:MET:SD	3:D:106:HIS:HE1	2.40	0.42
3:C:182:LEU:CD1	3:C:187:ILE:HB	2.50	0.42
3:C:49:GLU:HA	3:C:54:VAL:O	2.18	0.42
1:E:405:DG:C1'	1:E:406:DG:H5'	2.47	0.42
3:B:152:VAL:HG12	3:B:153:ILE:N	2.34	0.42
3:D:115:ARG:HB3	3:D:115:ARG:HE	1.63	0.42
3:D:29:ARG:NH1	3:D:32:GLU:OE2	2.52	0.42
3:A:136:ASP:O	3:A:139:GLY:N	2.52	0.42
3:B:17:GLU:O	3:B:20:GLU:HB2	2.19	0.42
3:B:92:LEU:HG	3:B:93:ASP:N	2.34	0.42
3:C:65:THR:HB	3:C:66:PRO:HD2	2.02	0.42
3:A:53:LEU:C	3:A:65:THR:HG23	2.40	0.42
3:C:110:ASP:O	3:C:113:GLU:HB2	2.19	0.42
3:C:77:ARG:HA	3:C:133:PRO:HD2	2.02	0.42
1:E:414:DA:C1'	1:E:415:DC:H5''	2.50	0.42
3:B:77:ARG:HA	3:B:133:PRO:HD2	2.01	0.42
3:D:11:TYR:OH	3:D:41:VAL:HA	2.20	0.42
3:D:59:ASP:O	3:D:60:ARG:HB2	2.19	0.41
3:C:57:ALA:HB2	3:C:63:GLN:NE2	2.34	0.41
3:A:107:VAL:HG21	3:B:107:VAL:HG21	2.02	0.41
3:A:166:VAL:O	3:A:167:GLN:HB3	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:196:VAL:O	3:B:202:ILE:HA	2.20	0.41
3:B:224:GLU:HG2	3:B:225:GLU:N	2.31	0.41
1:E:406:DG:H4'	3:A:60:ARG:NH2	2.35	0.41
3:B:163:VAL:HG22	3:B:164:ARG:N	2.35	0.41
3:B:178:GLN:OE1	3:B:214:LEU:HG	2.20	0.41
3:D:87:THR:CG2	3:D:97:VAL:HG21	2.47	0.41
3:C:213:LEU:CD1	3:C:218:ALA:HA	2.49	0.41
2:F:516:DC:H2''	3:D:40:THR:OG1	2.21	0.41
2:F:517:DC:C2'	2:F:518:DT:C7	2.98	0.41
3:B:182:LEU:HA	3:B:211:VAL:HG21	2.03	0.41
3:B:53:LEU:C	3:B:65:THR:HG23	2.41	0.41
3:C:29:ARG:HD3	3:C:29:ARG:HA	1.91	0.41
3:D:104:TRP:CD1	3:D:104:TRP:N	2.88	0.41
3:A:218:ALA:C	3:A:220:THR:H	2.24	0.41
3:B:110:ASP:O	3:B:113:GLU:HB2	2.20	0.41
3:B:13:ARG:NH1	3:B:17:GLU:OE1	2.52	0.41
3:B:201:HIS:N	3:B:201:HIS:CD2	2.89	0.41
3:C:86:LEU:HD22	3:C:92:LEU:CD2	2.46	0.41
3:D:5:VAL:HG11	3:D:106:HIS:HB2	2.03	0.41
3:A:17:GLU:O	3:A:20:GLU:HB2	2.20	0.41
3:B:214:LEU:HD12	3:B:217:LEU:CD2	2.51	0.41
2:F:506:DG:C2'	2:F:507:DT:H5''	2.47	0.41
3:B:194:GLU:O	3:B:204:LEU:HA	2.21	0.41
3:D:110:ASP:O	3:D:113:GLU:HB2	2.20	0.41
3:B:18:LEU:HD13	3:B:25:PRO:HA	2.03	0.41
3:B:54:VAL:CG2	3:B:55:VAL:N	2.82	0.41
3:C:89:ILE:N	3:C:89:ILE:HD12	2.35	0.41
3:D:85:LEU:HD22	3:D:116:LEU:HD21	2.03	0.41
3:B:95:ASN:N	3:B:95:ASN:OD1	2.54	0.41
3:A:29:ARG:HA	3:A:29:ARG:HD3	1.78	0.40
3:A:85:LEU:HD12	3:A:85:LEU:O	2.21	0.40
3:B:152:VAL:N	3:B:221:ILE:O	2.54	0.40
3:D:100:GLU:O	3:D:104:TRP:CD1	2.74	0.40
3:D:170:GLU:HA	3:D:173:GLN:CD	2.41	0.40
3:A:59:ASP:O	3:A:60:ARG:HB2	2.21	0.40
3:A:93:ASP:O	3:A:97:VAL:HG23	2.21	0.40
3:B:129:GLY:O	3:B:167:GLN:HG3	2.21	0.40
3:C:115:ARG:HD3	3:D:92:LEU:CA	2.47	0.40
3:A:78:LYS:HE3	3:A:108:MET:O	2.22	0.40
3:B:153:ILE:HB	3:B:218:ALA:HB1	2.04	0.40
3:C:90:ILE:HG12	3:C:90:ILE:H	1.72	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:401:DA:H2"	1:E:402:DT:H6	1.87	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	
3	A	167/225 (74%)	150 (90%)	14 (8%)	3 (2%)	8	41
3	B	215/225 (96%)	189 (88%)	21 (10%)	5 (2%)	6	34
3	C	169/225 (75%)	145 (86%)	21 (12%)	3 (2%)	8	41
3	D	169/225 (75%)	151 (89%)	14 (8%)	4 (2%)	6	34
All	All	720/900 (80%)	635 (88%)	70 (10%)	15 (2%)	7	37

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	121	LYS
3	B	121	LYS
3	C	121	LYS
3	D	121	LYS
3	A	3	ASP
3	B	3	ASP
3	C	3	ASP
3	D	3	ASP
3	D	175	GLU
3	A	175	GLU
3	B	175	GLU
3	B	176	THR
3	C	175	GLU
3	B	199	ASP

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Mol	Chain	Res	Type
3	D	169	ASN

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	
3	A	145/197 (74%)	112 (77%)	33 (23%)	1	4
3	B	188/197 (95%)	142 (76%)	46 (24%)	0	2
3	C	140/197 (71%)	112 (80%)	28 (20%)	1	6
3	D	139/197 (71%)	112 (81%)	27 (19%)	1	7
All	All	612/788 (78%)	478 (78%)	134 (22%)	1	5

All (134) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	13	ARG
3	A	21	GLU
3	A	27	ARG
3	A	29	ARG
3	A	33	ARG
3	A	37	SER
3	A	41	VAL
3	A	48	MET
3	A	67	THR
3	A	69	ARG
3	A	78	LYS
3	A	80	ARG
3	A	81	LEU
3	A	85	LEU
3	A	90	ILE
3	A	95	ASN
3	A	96	LYS
3	A	115	ARG
3	A	122	ASP
3	A	125	ARG

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Mol	Chain	Res	Type
3	A	132	ILE
3	A	137	GLU
3	A	140	VAL
3	A	165	ILE
3	A	166	VAL
3	A	169	ASN
3	A	171	ILE
3	A	176	THR
3	A	178	GLN
3	A	213	LEU
3	A	214	LEU
3	A	219	HIS
3	A	223	ILE
3	B	13	ARG
3	B	27	ARG
3	B	29	ARG
3	B	33	ARG
3	B	37	SER
3	B	41	VAL
3	B	48	MET
3	B	67	THR
3	B	69	ARG
3	B	78	LYS
3	B	80	ARG
3	B	81	LEU
3	B	85	LEU
3	B	90	ILE
3	B	95	ASN
3	B	96	LYS
3	B	115	ARG
3	B	122	ASP
3	B	125	ARG
3	B	128	PHE
3	B	132	ILE
3	B	137	GLU
3	B	140	VAL
3	B	152	VAL
3	B	157	THR
3	B	158	SER
3	B	161	ARG
3	B	167	GLN
3	B	171	ILE

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Mol	Chain	Res	Type
3	B	177	ASP
3	B	178	GLN
3	B	181	GLN
3	B	182	LEU
3	B	186	ASP
3	B	187	ILE
3	B	188	ARG
3	B	192	GLU
3	B	196	VAL
3	B	203	THR
3	B	204	LEU
3	B	205	SER
3	B	209	LYS
3	B	216	ASP
3	B	221	ILE
3	B	223	ILE
3	B	226	LEU
3	C	13	ARG
3	C	21	GLU
3	C	27	ARG
3	C	29	ARG
3	C	33	ARG
3	C	37	SER
3	C	41	VAL
3	C	48	MET
3	C	67	THR
3	C	69	ARG
3	C	78	LYS
3	C	80	ARG
3	C	81	LEU
3	C	85	LEU
3	C	90	ILE
3	C	95	ASN
3	C	96	LYS
3	C	115	ARG
3	C	122	ASP
3	C	125	ARG
3	C	128	PHE
3	C	132	ILE
3	C	169	ASN
3	C	171	ILE
3	C	176	THR

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Mol	Chain	Res	Type
3	C	178	GLN
3	C	213	LEU
3	C	214	LEU
3	D	13	ARG
3	D	21	GLU
3	D	27	ARG
3	D	29	ARG
3	D	33	ARG
3	D	37	SER
3	D	48	MET
3	D	67	THR
3	D	69	ARG
3	D	78	LYS
3	D	80	ARG
3	D	81	LEU
3	D	85	LEU
3	D	90	ILE
3	D	95	ASN
3	D	96	LYS
3	D	115	ARG
3	D	122	ASP
3	D	124	SER
3	D	125	ARG
3	D	132	ILE
3	D	169	ASN
3	D	171	ILE
3	D	176	THR
3	D	178	GLN
3	D	213	LEU
3	D	214	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	B	181	GLN
3	B	201	HIS
3	B	207	ASN

5.3.3 RNA ⓘ

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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	E	21/21 (100%)	0.28	0 100 100	21, 33, 45, 49	2 (9%)
2	F	21/21 (100%)	0.31	0 100 100	23, 36, 41, 49	2 (9%)
3	A	173/225 (76%)	0.10	14 (8%) 12 6	16, 44, 93, 100	0
3	B	219/225 (97%)	-0.16	2 (0%) 84 75	14, 42, 84, 100	0
3	C	175/225 (77%)	0.04	6 (3%) 45 29	13, 59, 94, 100	0
3	D	175/225 (77%)	0.58	14 (8%) 12 6	21, 61, 100, 100	0
All	All	784/942 (83%)	0.13	36 (4%) 32 20	13, 50, 95, 100	4 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	167	GLN	7.1
3	D	220	THR	4.3
3	A	220	THR	4.0
3	D	168	ILE	4.0
3	D	187	ILE	4.0
3	D	221	ILE	3.9
3	D	213	LEU	3.8
3	A	168	ILE	3.6
3	A	169	ASN	3.6
3	D	182	LEU	3.4
3	A	218	ALA	3.3
3	A	213	LEU	3.3
3	D	169	ASN	3.2
3	A	217	LEU	3.1
3	C	167	GLN	2.9
3	D	166	VAL	2.9
3	C	221	ILE	2.9
3	A	182	LEU	2.7
3	D	172	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
3	A	178	GLN	2.5
3	B	199	ASP	2.5
3	D	217	LEU	2.5
3	D	93	ASP	2.4
3	D	218	ALA	2.4
3	A	167	GLN	2.4
3	A	180	THR	2.3
3	A	179	PHE	2.3
3	C	179	PHE	2.3
3	B	198	ARG	2.3
3	A	93	ASP	2.2
3	D	16	TYR	2.2
3	C	178	GLN	2.2
3	C	168	ILE	2.1
3	A	211	VAL	2.1
3	A	181	GLN	2.0
3	C	182	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
4	CO	B	602	1/1	0.96	0.15	37,37,37,37	0
4	CO	A	601	1/1	0.97	0.07	38,38,38,38	0
4	CO	C	603	1/1	0.97	0.09	50,50,50,50	0
4	CO	D	604	1/1	0.98	0.02	54,54,54,54	0
4	CO	C	703	1/1	0.98	0.18	50,50,50,50	0
4	CO	A	701	1/1	0.98	0.17	38,38,38,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CO	D	704	1/1	0.99	0.10	54,54,54,54	0
4	CO	B	702	1/1	0.99	0.15	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.