



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:09 PM UTC

PDB ID : 3RN5 / pdb_00003rn5
Title : Structural basis of cytosolic DNA recognition by innate immune receptors
Authors : Jin, T.C.; Xiao, T.
Deposited on : 2011-04-21
Resolution : 2.50 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

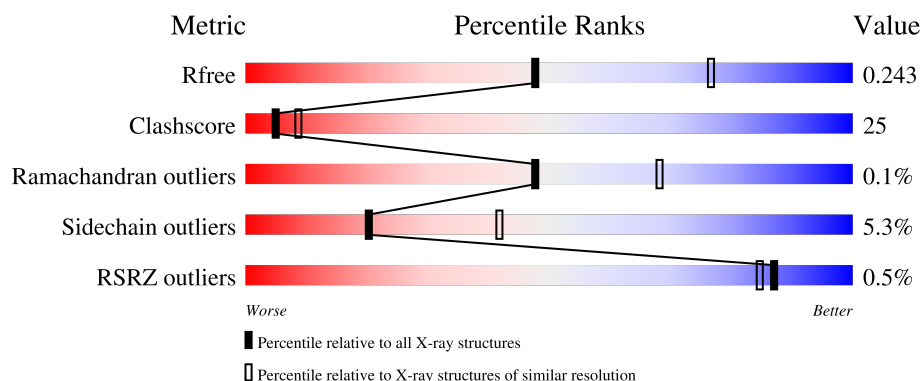
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

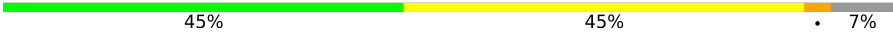
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	208	
1	B	208	
1	C	208	
1	D	208	
2	K	19	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	M	19	
3	L	19	
3	N	19	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	EDO	D	6	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7944 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 Interferon-inducible protein AIM2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	194	Total	C	N	O	S	0	0	0
			1552	998	268	280	6			
1	B	194	Total	C	N	O	S	0	0	0
			1552	998	268	280	6			
1	C	194	Total	C	N	O	S	0	0	0
			1552	998	268	280	6			
1	D	193	Total	C	N	O	S	0	0	0
			1543	993	267	277	6			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	GLY	-	expression tag	UNP O14862
A	141	SER	-	expression tag	UNP O14862
A	142	VAL	-	expression tag	UNP O14862
A	143	ASP	-	expression tag	UNP O14862
A	344	ALA	-	expression tag	UNP O14862
A	345	ALA	-	expression tag	UNP O14862
A	346	ALA	-	expression tag	UNP O14862
A	347	SER	-	expression tag	UNP O14862
B	140	GLY	-	expression tag	UNP O14862
B	141	SER	-	expression tag	UNP O14862
B	142	VAL	-	expression tag	UNP O14862
B	143	ASP	-	expression tag	UNP O14862
B	344	ALA	-	expression tag	UNP O14862
B	345	ALA	-	expression tag	UNP O14862
B	346	ALA	-	expression tag	UNP O14862
B	347	SER	-	expression tag	UNP O14862
C	140	GLY	-	expression tag	UNP O14862
C	141	SER	-	expression tag	UNP O14862
C	142	VAL	-	expression tag	UNP O14862
C	143	ASP	-	expression tag	UNP O14862
C	344	ALA	-	expression tag	UNP O14862

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	345	ALA	-	expression tag	UNP O14862
C	346	ALA	-	expression tag	UNP O14862
C	347	SER	-	expression tag	UNP O14862
D	140	GLY	-	expression tag	UNP O14862
D	141	SER	-	expression tag	UNP O14862
D	142	VAL	-	expression tag	UNP O14862
D	143	ASP	-	expression tag	UNP O14862
D	344	ALA	-	expression tag	UNP O14862
D	345	ALA	-	expression tag	UNP O14862
D	346	ALA	-	expression tag	UNP O14862
D	347	SER	-	expression tag	UNP O14862

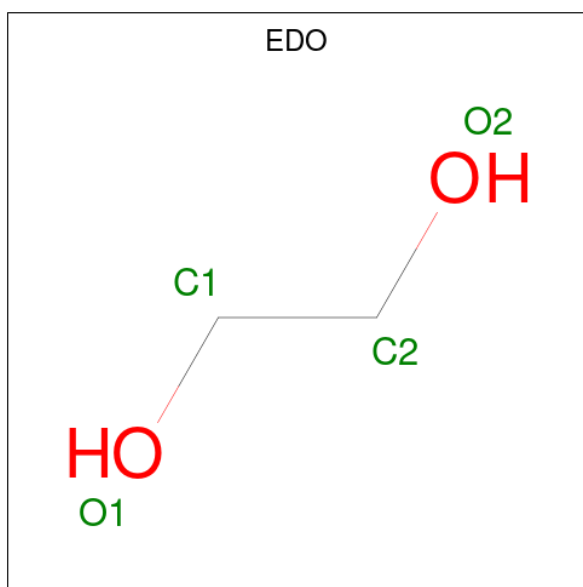
- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*AP*TP*CP*AP*AP*AP*GP*AP*GP*AP*GP*AP*GP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	19	Total	C	N	O	P	0	0	0
			394	187	86	103	18			
2	M	19	Total	C	N	O	P	0	0	0
			394	187	86	103	18			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	19	Total	C	N	O	P	0	0	0
			379	185	55	121	18			
3	N	19	Total	C	N	O	P	0	0	0
			379	185	55	121	18			

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	37	Total	O	0	0
			37	37		
5	B	39	Total	O	0	0
			39	39		
5	C	27	Total	O	0	0
			27	27		

Continued on next page...

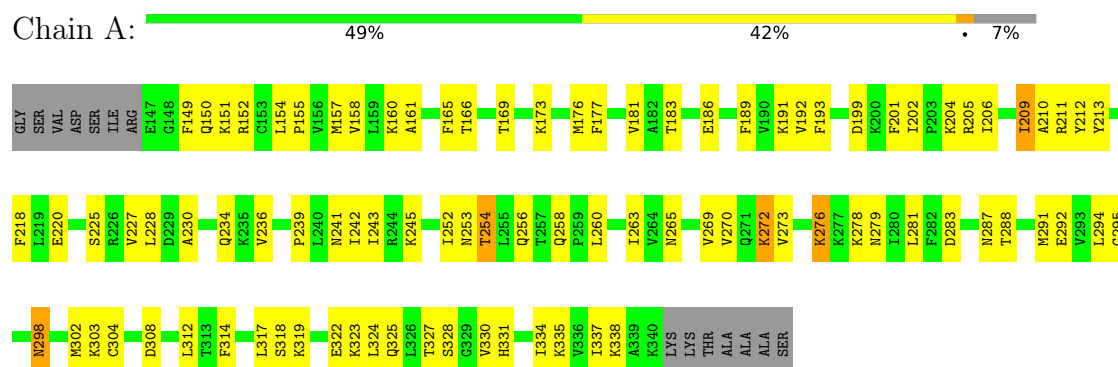
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	25	Total 25	O 25	0	0
5	K	14	Total 14	O 14	0	0
5	L	7	Total 7	O 7	0	0
5	M	5	Total 5	O 5	0	0
5	N	9	Total 9	O 9	0	0

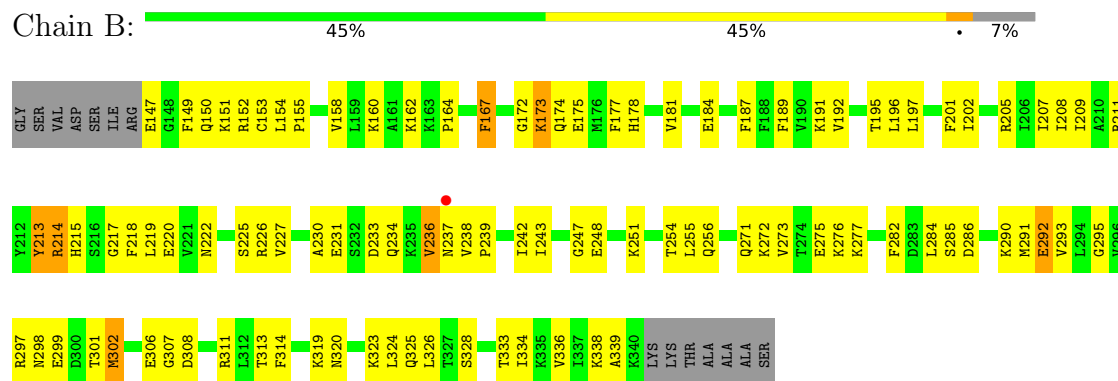
3 Residue-property plots

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.

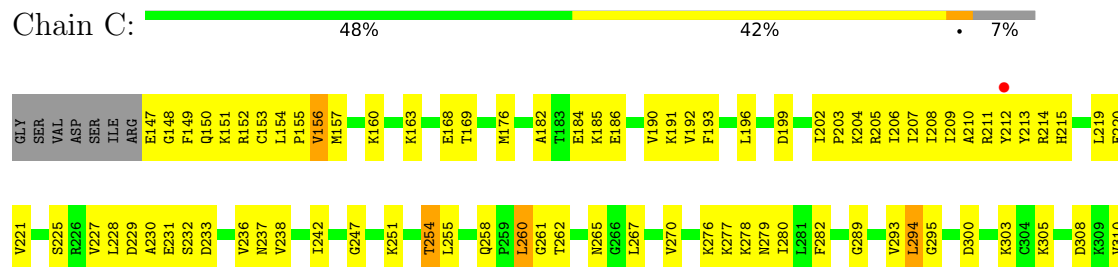
• Molecule 1: Interferon-inducible protein AIM2



• Molecule 1: Interferon-inducible protein AIM2



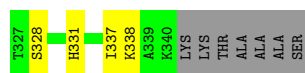
• Molecule 1: Interferon-inducible protein AIM2





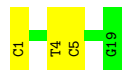
- Molecule 1: Interferon-inducible protein AIM2

Chain D: 54% 38% 7%



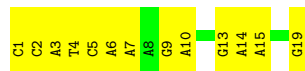
- Molecule 2: DNA (5'-D(*CP*CP*AP*TP*CP*AP*AP*AP*GP*AP*GP*AP*GP*AP*AP*AP*GP*AP*G)-3')

Chain K: 84% 16%



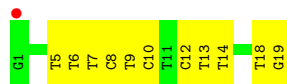
- Molecule 2: DNA (5'-D(*CP*CP*AP*TP*CP*AP*AP*AP*GP*AP*GP*AP*GP*AP*AP*AP*GP*AP*G)-3')

Chain M: 32% 68%



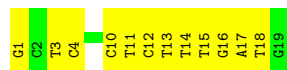
- Molecule 3: DNA (5'-D(*GP*CP*TP*CP*TP*TP*TP*CP*TP*CP*TP*CP*TP*TP*TP*GP*AP*TP*G)-3')

Chain L: 5% 42% 58%



- Molecule 3: DNA (5'-D(*GP*CP*TP*CP*TP*TP*TP*CP*TP*CP*TP*CP*TP*TP*TP*GP*AP*TP*G)-3')

Chain N: 37% 63%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.36Å 136.36Å 77.04Å 90.00° 89.98° 90.00°	Depositor
Resolution (Å)	46.02 – 2.50 46.02 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.02-2.50) 94.1 (46.02-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_736)	Depositor
R, R_{free}	0.201 , 0.241 0.201 , 0.243	Depositor DCC
R_{free} test set	2234 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.645	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 61.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.449 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7944	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.0682e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: EDO

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	0.40	0/1576	0.82	2/2113 (0.1%)
1	B	0.41	0/1576	0.81	2/2113 (0.1%)
1	C	0.40	0/1576	0.84	2/2113 (0.1%)
1	D	0.40	0/1567	0.81	2/2101 (0.1%)
2	K	0.13	0/446	0.54	0/687
2	M	0.11	0/446	0.49	0/687
3	L	0.13	0/420	0.55	0/646
3	N	0.13	0/420	0.56	0/646
All	All	0.36	0/8027	0.76	8/11106 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	289	GLY	N-CA-C	6.46	118.84	110.45
1	B	213	TYR	N-CA-C	6.41	120.31	112.23
1	C	289	GLY	N-CA-C	6.09	118.18	110.38
1	C	261	GLY	N-CA-C	-5.54	108.39	115.42
1	A	298	ASN	N-CA-C	5.44	115.42	108.24
1	B	214	ARG	N-CA-C	5.38	117.20	110.91
1	A	186	GLU	N-CA-C	5.22	116.92	109.14
1	D	232	SER	N-CA-C	-5.19	106.64	112.87

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	1552	0	1636	88	0
1	B	1552	0	1636	94	0
1	C	1552	0	1636	92	0
1	D	1543	0	1630	80	0
2	K	394	0	212	2	0
2	M	394	0	212	16	0
3	L	379	0	221	11	0
3	N	379	0	221	13	0
4	A	8	0	12	4	0
4	B	8	0	12	3	0
4	C	4	0	6	0	0
4	D	16	0	24	9	0
5	A	37	0	0	4	0
5	B	39	0	0	4	0
5	C	27	0	0	4	0
5	D	25	0	0	2	0
5	K	14	0	0	1	0
5	L	7	0	0	0	0
5	M	5	0	0	0	0
5	N	9	0	0	1	0
All	All	7944	0	7458	382	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 25.

All (382) 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:B:291:MET:SD	1:B:324:LEU:HD11	1.77	1.24
1:A:230:ALA:HB1	1:A:234:GLN:HG3	1.44	0.98
1:B:230:ALA:HB1	1:B:234:GLN:HG3	1.48	0.95
1:D:262:THR:HA	4:D:6:EDO:H21	1.49	0.94
1:A:158:VAL:HA	1:A:181:VAL:HG12	1.52	0.92
1:A:283:ASP:OD1	1:A:292:GLU:HG2	1.69	0.92
1:B:158:VAL:HA	1:B:181:VAL:HG12	1.54	0.88

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:MET:SD	1:B:324:LEU:CD1	2.62	0.87
1:D:157:MET:HE3	1:D:204:LYS:HA	1.57	0.87
3:L:5:DT:H2'	3:L:6:DT:H72	1.57	0.85
1:A:323:LYS:O	1:A:323:LYS:HG3	1.77	0.84
1:B:155:PRO:HB2	1:B:236:VAL:HG21	1.62	0.81
1:C:319:LYS:HB3	1:C:325:GLN:HB3	1.62	0.81
1:C:276:LYS:HE3	1:C:294:LEU:HD11	1.64	0.79
1:C:157:MET:HE3	1:C:204:LYS:HA	1.64	0.78
1:C:251:LYS:O	1:C:254:THR:HB	1.83	0.78
1:A:322:GLU:O	1:A:323:LYS:HG2	1.83	0.78
1:B:208:ILE:HG13	1:B:230:ALA:HB2	1.65	0.78
1:B:255:LEU:HB2	1:B:324:LEU:CD2	2.13	0.77
2:M:6:DA:H2''	2:M:7:DA:H5'	1.64	0.77
1:C:192:VAL:HA	1:C:221:VAL:HB	1.66	0.77
1:C:147:GLU:N	1:C:214:ARG:HE	1.82	0.76
1:D:296:VAL:HG12	1:D:297:ARG:H	1.49	0.76
1:C:154:LEU:HD12	1:C:155:PRO:HD2	1.68	0.76
1:D:240:LEU:HD12	1:D:240:LEU:H	1.51	0.76
1:B:239:PRO:HD2	1:B:242:ILE:HD12	1.68	0.75
1:C:255:LEU:HA	1:C:258:GLN:HG2	1.68	0.75
1:B:202:ILE:HD12	1:B:205:ARG:HD3	1.68	0.74
1:A:211:ARG:HB2	1:A:225:SER:HA	1.70	0.74
1:A:160:LYS:HE3	4:A:7:EDO:H11	1.70	0.74
1:A:209:ILE:HG23	1:A:227:VAL:HG22	1.71	0.73
1:A:294:LEU:HB2	5:A:122:HOH:O	1.88	0.73
3:N:16:DG:H2''	3:N:17:DA:H5'	1.72	0.72
1:B:255:LEU:HB2	1:B:324:LEU:HD21	1.71	0.71
1:B:255:LEU:CB	1:B:324:LEU:HD22	2.20	0.71
3:L:8:DC:H4'	3:L:9:DT:OP1	1.90	0.71
1:A:241:ASN:O	1:A:245:LYS:HG2	1.91	0.71
1:D:315:PHE:CE2	1:D:328:SER:HB3	2.25	0.70
1:D:324:LEU:HD23	1:D:324:LEU:H	1.54	0.70
1:A:287:ASN:HD22	1:D:240:LEU:HD13	1.56	0.70
1:D:313:THR:HG23	4:D:9:EDO:H21	1.72	0.70
1:A:158:VAL:HG11	1:A:202:ILE:H	1.57	0.70
1:B:211:ARG:HE	1:B:226:ARG:HD2	1.54	0.70
1:B:324:LEU:C	1:B:324:LEU:HD12	2.17	0.69
1:C:151:LYS:HG2	5:C:84:HOH:O	1.93	0.69
1:C:319:LYS:HB3	1:C:325:GLN:N	2.06	0.69
1:D:149:PHE:HB3	1:D:213:TYR:CD1	2.28	0.69
1:B:291:MET:HB2	1:B:324:LEU:O	1.94	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:5:DC:H1'	2:M:6:DA:C8	2.28	0.68
1:D:231:GLU:CD	1:D:231:GLU:H	2.01	0.68
1:C:151:LYS:CG	5:C:84:HOH:O	2.41	0.68
1:B:292:GLU:O	1:B:325:GLN:HG3	1.93	0.68
1:D:149:PHE:HB2	1:D:213:TYR:HA	1.76	0.67
1:D:149:PHE:HB3	1:D:213:TYR:CG	2.30	0.67
3:L:5:DT:H2'	3:L:6:DT:C7	2.25	0.66
3:L:9:DT:H1'	3:L:10:DC:H5''	1.77	0.66
3:N:11:DT:H1'	3:N:12:DC:H5''	1.76	0.66
1:D:309:LYS:HE2	5:N:150:HOH:O	1.96	0.66
1:B:151:LYS:HE3	1:B:211:ARG:NH1	2.12	0.65
1:C:211:ARG:HB2	1:C:225:SER:HA	1.79	0.65
1:B:297:ARG:NE	1:B:302:MET:HE1	2.12	0.64
1:C:319:LYS:CB	1:C:325:GLN:HB3	2.26	0.64
1:B:255:LEU:HB3	1:B:324:LEU:HD22	1.79	0.63
1:C:255:LEU:HA	1:C:258:GLN:CG	2.28	0.63
1:D:296:VAL:HG12	1:D:297:ARG:N	2.14	0.63
1:D:325:GLN:HG2	1:D:326:LEU:N	2.13	0.63
1:C:186:GLU:HA	1:C:242:ILE:HD13	1.80	0.62
1:D:149:PHE:HB2	1:D:212:TYR:O	1.99	0.62
1:C:148:GLY:O	1:C:214:ARG:HB3	2.00	0.62
1:A:201:PHE:C	1:A:202:ILE:HG13	2.25	0.62
1:C:319:LYS:HB3	1:C:325:GLN:H	1.65	0.62
1:B:255:LEU:CB	1:B:324:LEU:CD2	2.77	0.61
1:A:252:ILE:HG23	1:A:324:LEU:HD21	1.82	0.61
1:A:166:THR:OG1	1:A:173:LYS:HE2	2.00	0.61
1:B:175:GLU:HB2	1:B:195:THR:OG1	2.00	0.61
1:A:303:LYS:HE2	1:A:338:LYS:NZ	2.15	0.61
1:A:324:LEU:H	1:A:324:LEU:HD23	1.64	0.61
1:D:176:MET:SD	1:D:331:HIS:ND1	2.74	0.61
1:C:305:LYS:HG2	1:C:308:ASP:OD2	2.01	0.60
1:D:149:PHE:CB	1:D:213:TYR:HA	2.31	0.60
1:B:207:ILE:HD11	1:B:209:ILE:HD11	1.83	0.60
1:D:282:PHE:HE2	1:D:295:GLY:HA3	1.66	0.60
1:C:247:GLY:HA3	2:M:15:DA:OP1	2.02	0.60
1:D:262:THR:HG23	4:D:6:EDO:H12	1.83	0.60
1:A:158:VAL:HG11	1:A:202:ILE:N	2.15	0.59
1:B:205:ARG:HH21	1:B:207:ILE:HG22	1.67	0.59
1:A:150:GLN:HG3	1:A:212:TYR:CZ	2.38	0.59
1:C:168:GLU:O	1:C:169:THR:HG23	2.03	0.59
1:A:323:LYS:O	1:A:323:LYS:CG	2.47	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:1:DC:H2''	2:M:2:DC:C5'	2.32	0.59
2:M:6:DA:C2'	2:M:7:DA:H5'	2.32	0.59
3:L:5:DT:H2''	3:L:6:DT:C6	2.38	0.58
1:A:210:ALA:O	1:A:225:SER:HB2	2.03	0.58
1:B:151:LYS:HE3	1:B:211:ARG:HH12	1.66	0.58
1:C:152:ARG:HG2	1:C:153:CYS:H	1.68	0.58
1:D:193:PHE:HE2	1:D:220:GLU:HG3	1.67	0.58
1:D:251:LYS:O	1:D:254:THR:HB	2.04	0.57
1:A:150:GLN:HG3	1:A:212:TYR:CE1	2.38	0.57
1:C:193:PHE:HE2	1:C:220:GLU:HG3	1.67	0.57
1:C:260:LEU:HD22	1:C:318:SER:O	2.04	0.57
1:C:260:LEU:N	1:C:260:LEU:HD23	2.18	0.57
1:D:214:ARG:HH21	1:D:218:PHE:C	2.12	0.57
1:D:296:VAL:HG23	5:D:87:HOH:O	2.04	0.57
1:C:184:GLU:HG2	1:C:237:ASN:O	2.04	0.57
1:D:298:ASN:O	1:D:302:MET:HG3	2.05	0.56
1:C:163:LYS:HG2	3:N:10:DC:H3'	1.86	0.56
1:C:202:ILE:HB	1:C:205:ARG:HG3	1.87	0.56
1:C:300:ASP:HA	1:C:303:LYS:HD3	1.87	0.56
2:M:1:DC:H2''	2:M:2:DC:H5'	1.87	0.56
1:A:256:GLN:HG2	1:A:324:LEU:HD22	1.88	0.56
1:C:148:GLY:O	1:C:214:ARG:HD3	2.05	0.56
1:D:211:ARG:HB2	1:D:225:SER:HA	1.88	0.56
1:D:252:ILE:HG13	1:D:286:ASP:OD2	2.06	0.56
1:A:327:THR:HG22	5:A:122:HOH:O	2.06	0.56
1:D:262:THR:HA	4:D:6:EDO:C2	2.32	0.56
1:A:260:LEU:HD23	1:A:317:LEU:HG	1.88	0.55
1:C:206:ILE:HG22	1:C:230:ALA:HB2	1.86	0.55
1:A:287:ASN:ND2	1:D:240:LEU:HD13	2.21	0.55
1:C:196:LEU:HD22	1:C:196:LEU:H	1.72	0.55
1:A:230:ALA:HB1	1:A:234:GLN:CG	2.29	0.55
1:B:307:GLY:O	1:B:339:ALA:HB2	2.07	0.55
1:C:196:LEU:HD22	1:C:196:LEU:N	2.22	0.55
1:B:164:PRO:HA	1:B:177:PHE:HB3	1.87	0.55
2:M:5:DC:H4'	2:M:6:DA:OP1	2.07	0.55
1:C:280:ILE:N	1:C:295:GLY:O	2.40	0.55
1:D:338:LYS:HE2	1:D:338:LYS:HA	1.89	0.55
1:C:149:PHE:HD1	1:C:213:TYR:CG	2.25	0.54
1:B:154:LEU:HD12	1:B:155:PRO:HD2	1.88	0.54
1:B:214:ARG:HA	1:B:219:LEU:HD22	1.89	0.54
1:C:202:ILE:O	1:C:205:ARG:HB2	2.07	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ASP:HB2	5:A:38:HOH:O	2.07	0.54
1:B:162:LYS:HD2	4:B:8:EDO:H11	1.89	0.54
1:D:231:GLU:C	1:D:233:ASP:N	2.64	0.54
1:B:208:ILE:CG1	1:B:230:ALA:HB2	2.36	0.54
1:C:319:LYS:HG3	1:C:324:LEU:HA	1.89	0.54
1:C:255:LEU:HD23	1:C:258:GLN:HG3	1.89	0.53
1:A:295:GLY:HA2	1:A:328:SER:HB2	1.88	0.53
1:C:317:LEU:HD12	1:C:326:LEU:CD2	2.39	0.53
2:M:9:DG:H2''	2:M:10:DA:OP2	2.09	0.53
3:N:17:DA:H1'	3:N:18:DT:H5'	1.89	0.53
1:A:292:GLU:O	1:A:325:GLN:HG3	2.07	0.53
1:B:217:GLY:HA2	5:B:348:HOH:O	2.07	0.53
1:C:156:VAL:HA	1:C:236:VAL:HG11	1.90	0.53
1:A:241:ASN:HD21	1:D:253:ASN:ND2	2.06	0.53
1:A:241:ASN:HD21	1:D:253:ASN:HD21	1.57	0.53
1:A:281:LEU:HD11	1:A:292:GLU:HB3	1.91	0.53
1:A:151:LYS:O	1:A:152:ARG:HG3	2.08	0.53
1:A:201:PHE:O	1:A:202:ILE:HG13	2.09	0.53
1:B:192:VAL:HG11	1:B:201:PHE:CE2	2.43	0.53
1:D:187:PHE:HE1	1:D:245:LYS:HE3	1.74	0.53
1:A:239:PRO:HG2	1:A:242:ILE:HG13	1.90	0.52
1:B:191:LYS:HE3	1:B:314:PHE:CE1	2.43	0.52
1:D:231:GLU:C	1:D:233:ASP:H	2.17	0.52
1:A:331:HIS:HB2	4:A:2:EDO:O2	2.09	0.52
1:B:211:ARG:NE	1:B:226:ARG:HD2	2.23	0.52
1:A:241:ASN:HD21	1:D:253:ASN:CG	2.18	0.52
1:C:149:PHE:HE2	1:C:151:LYS:HB3	1.75	0.52
1:C:152:ARG:O	1:C:210:ALA:HA	2.10	0.52
1:A:157:MET:HE3	1:A:204:LYS:HA	1.91	0.52
1:A:158:VAL:HB	1:A:202:ILE:O	2.10	0.52
1:B:178:HIS:HE1	1:B:333:THR:HG23	1.75	0.52
2:M:5:DC:H1'	2:M:6:DA:N7	2.24	0.52
1:C:319:LYS:HB3	1:C:325:GLN:CB	2.38	0.51
1:C:185:LYS:O	1:C:186:GLU:HB3	2.10	0.51
1:D:217:GLY:CA	4:D:6:EDO:H22	2.41	0.51
1:C:317:LEU:HD12	1:C:326:LEU:HD21	1.92	0.51
1:C:203:PRO:O	1:C:204:LYS:HB2	2.10	0.51
1:B:150:GLN:O	1:B:211:ARG:HA	2.10	0.51
1:B:184:GLU:HG2	1:B:237:ASN:N	2.26	0.51
1:D:282:PHE:CE2	1:D:295:GLY:HA3	2.45	0.51
1:A:160:LYS:HG2	4:A:7:EDO:H11	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ASN:ND2	1:D:253:ASN:HD21	2.10	0.50
1:C:151:LYS:HG3	5:C:84:HOH:O	2.08	0.50
1:A:206:ILE:HG22	1:A:230:ALA:HB3	1.94	0.50
1:B:197:LEU:HB3	1:B:201:PHE:HE2	1.77	0.50
1:D:193:PHE:CE2	1:D:220:GLU:HG3	2.46	0.50
1:B:319:LYS:HB3	1:B:325:GLN:H	1.77	0.50
1:C:228:LEU:HG	1:C:229:ASP:H	1.75	0.50
1:C:209:ILE:HG12	1:C:227:VAL:HG22	1.93	0.49
1:C:152:ARG:HG2	1:C:153:CYS:N	2.27	0.49
1:C:150:GLN:HB2	1:C:214:ARG:HD2	1.93	0.49
3:N:14:DT:H1'	3:N:15:DT:H5''	1.95	0.49
1:C:154:LEU:HD12	1:C:155:PRO:CD	2.42	0.49
1:D:296:VAL:CG1	1:D:297:ARG:H	2.23	0.49
1:B:256:GLN:O	1:B:320:ASN:HB2	2.12	0.49
1:A:211:ARG:O	1:A:225:SER:HB3	2.11	0.49
1:C:163:LYS:HG3	3:N:10:DC:P	2.53	0.49
1:D:166:THR:HG22	1:D:175:GLU:HB3	1.94	0.49
1:D:177:PHE:CE2	1:D:192:VAL:HB	2.47	0.49
1:A:270:VAL:HG21	1:A:304:CYS:SG	2.53	0.49
1:D:194:ASN:O	1:D:197:LEU:HB2	2.13	0.49
1:D:221:VAL:HG13	1:D:225:SER:OG	2.12	0.49
1:D:187:PHE:C	1:D:187:PHE:CD2	2.91	0.48
1:B:187:PHE:C	1:B:187:PHE:CD2	2.91	0.48
1:B:189:PHE:O	1:B:218:PHE:HB2	2.14	0.48
1:B:290:LYS:HE3	1:B:323:LYS:HE2	1.95	0.48
1:B:290:LYS:CE	1:B:323:LYS:HE2	2.44	0.48
1:C:315:PHE:CE2	1:C:328:SER:HB2	2.48	0.48
1:C:230:ALA:C	1:C:232:SER:H	2.20	0.48
1:D:184:GLU:HG2	1:D:237:ASN:O	2.14	0.47
3:N:13:DT:H2'	3:N:14:DT:H72	1.94	0.47
1:A:213:TYR:HB3	1:A:220:GLU:HB3	1.96	0.47
1:B:215:HIS:HE2	1:B:220:GLU:CB	2.27	0.47
1:C:156:VAL:C	1:C:236:VAL:HG11	2.39	0.47
1:A:169:THR:HG22	5:A:349:HOH:O	2.14	0.47
1:B:222:ASN:O	1:B:225:SER:HB3	2.13	0.47
1:C:149:PHE:CE2	1:C:151:LYS:HB3	2.49	0.47
1:C:193:PHE:HD2	1:C:221:VAL:O	1.98	0.47
1:B:231:GLU:C	1:B:233:ASP:N	2.71	0.47
1:C:207:ILE:C	1:C:208:ILE:HD13	2.38	0.47
1:D:217:GLY:HA2	4:D:6:EDO:H22	1.97	0.47
1:D:234:GLN:OE1	1:D:235:LYS:N	2.48	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:LEU:HD12	1:A:155:PRO:HD2	1.97	0.47
1:C:156:VAL:CA	1:C:236:VAL:HG11	2.44	0.47
2:M:3:DA:H2''	2:M:4:DT:O5'	2.15	0.47
2:M:9:DG:H1'	2:M:10:DA:H5'	1.97	0.47
1:A:191:LYS:HD2	1:A:218:PHE:CD1	2.50	0.47
1:C:190:VAL:HG22	1:C:219:LEU:HB2	1.97	0.47
1:D:274:THR:O	1:D:281:LEU:N	2.48	0.47
1:A:160:LYS:HE3	4:A:7:EDO:C1	2.44	0.46
1:C:176:MET:HE1	1:C:191:LYS:HB3	1.97	0.46
1:C:267:LEU:HD12	1:C:310:VAL:O	2.14	0.46
1:A:303:LYS:HB3	1:A:338:LYS:CE	2.45	0.46
1:B:207:ILE:HG13	1:B:209:ILE:HG13	1.97	0.46
1:D:279:ASN:HB3	1:D:294:LEU:HG	1.98	0.46
1:A:303:LYS:HE2	1:A:338:LYS:HZ1	1.80	0.46
1:A:308:ASP:OD2	1:A:338:LYS:HD3	2.16	0.46
1:B:160:LYS:HD2	4:B:8:EDO:C1	2.45	0.46
1:B:173:LYS:O	1:B:173:LYS:HG3	2.16	0.46
1:D:298:ASN:OD1	1:D:301:THR:N	2.35	0.46
1:A:160:LYS:HG3	1:A:161:ALA:N	2.30	0.46
1:B:319:LYS:HG2	1:B:320:ASN:N	2.31	0.46
1:B:319:LYS:O	1:B:324:LEU:HB2	2.16	0.46
1:D:149:PHE:CD2	1:D:213:TYR:CZ	3.04	0.46
2:K:4:DT:H2''	2:K:5:DC:H5'	1.98	0.46
1:B:308:ASP:CG	1:B:338:LYS:HD3	2.40	0.45
1:A:149:PHE:CZ	1:A:211:ARG:HA	2.51	0.45
1:B:247:GLY:HA3	3:L:18:DT:OP1	2.17	0.45
1:B:293:VAL:HA	1:B:326:LEU:O	2.16	0.45
1:C:193:PHE:CE2	1:C:220:GLU:HG3	2.50	0.45
1:C:228:LEU:HG	1:C:229:ASP:N	2.31	0.45
1:C:270:VAL:HB	1:C:308:ASP:O	2.17	0.45
1:A:298:ASN:O	1:A:302:MET:HG3	2.16	0.45
1:A:206:ILE:HG22	1:A:230:ALA:CB	2.46	0.45
1:B:256:GLN:HA	1:B:320:ASN:HB2	1.98	0.45
1:C:260:LEU:C	1:C:262:THR:H	2.25	0.45
2:K:1:DC:H5''	2:M:19:DG:H2''	1.98	0.45
1:A:155:PRO:HB2	1:A:236:VAL:CG2	2.47	0.45
1:B:319:LYS:HB3	1:B:325:GLN:N	2.32	0.45
1:A:254:THR:O	1:A:258:GLN:HG2	2.16	0.45
1:B:158:VAL:HB	1:B:202:ILE:O	2.16	0.45
1:B:209:ILE:HG23	1:B:227:VAL:HG22	1.98	0.45
1:B:238:VAL:HG12	1:B:243:ILE:HG13	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:TYR:CD2	1:C:212:TYR:N	2.84	0.45
1:C:231:GLU:O	1:C:232:SER:OG	2.28	0.45
1:C:231:GLU:C	1:C:233:ASP:H	2.24	0.45
1:D:265:ASN:HB3	1:D:313:THR:HA	1.98	0.44
1:A:151:LYS:HA	1:A:211:ARG:H	1.82	0.44
1:A:263:ILE:HG23	1:A:314:PHE:HA	1.99	0.44
1:B:295:GLY:HA2	1:B:328:SER:HB2	1.99	0.44
1:D:231:GLU:HB2	1:D:233:ASP:HB3	1.99	0.44
3:N:10:DC:H2'	3:N:11:DT:H72	1.97	0.44
1:A:158:VAL:HG23	1:A:205:ARG:O	2.18	0.44
1:B:218:PHE:O	1:B:219:LEU:HD23	2.16	0.44
1:A:256:GLN:CG	1:A:324:LEU:HD22	2.48	0.44
1:A:204:LYS:HE2	3:L:12:DC:OP1	2.17	0.44
1:A:260:LEU:CD2	1:A:317:LEU:HG	2.48	0.44
1:A:278:LYS:HD3	1:A:278:LYS:HA	1.76	0.44
1:B:152:ARG:H	1:B:152:ARG:CD	2.31	0.44
1:D:291:MET:SD	1:D:326:LEU:HD11	2.57	0.44
3:L:13:DT:H2''	3:L:14:DT:H5'	1.98	0.44
1:A:265:ASN:HA	1:A:312:LEU:O	2.18	0.44
1:B:196:LEU:N	1:B:196:LEU:HD22	2.33	0.44
1:A:189:PHE:CE2	1:A:265:ASN:ND2	2.86	0.43
1:A:335:LYS:HD3	5:K:64:HOH:O	2.17	0.43
1:C:230:ALA:C	1:C:232:SER:N	2.75	0.43
1:D:210:ALA:O	1:D:225:SER:HB2	2.18	0.43
1:A:227:VAL:O	1:A:228:LEU:HG	2.18	0.43
1:B:282:PHE:HB3	1:B:284:LEU:HD21	1.99	0.43
1:D:313:THR:CG2	4:D:9:EDO:H21	2.45	0.43
1:A:260:LEU:HD13	1:A:318:SER:O	2.19	0.43
1:B:214:ARG:HG3	1:B:219:LEU:HD22	2.00	0.43
1:C:214:ARG:O	1:C:214:ARG:HG2	2.19	0.43
3:L:7:DT:H1'	3:L:8:DC:H5'	1.99	0.43
1:B:167:PHE:N	1:B:167:PHE:CD2	2.87	0.43
1:D:160:LYS:HD3	5:D:123:HOH:O	2.17	0.43
1:B:172:GLY:O	1:B:174:GLN:HG2	2.17	0.43
1:C:149:PHE:HA	1:C:213:TYR:HA	2.00	0.43
1:D:276:LYS:HG3	1:D:294:LEU:HD11	2.00	0.43
1:A:165:PHE:CZ	1:A:176:MET:SD	3.11	0.43
1:B:301:THR:HA	1:B:336:VAL:HG21	2.00	0.43
3:N:3:DT:H1'	3:N:4:DC:H5''	1.99	0.43
1:A:191:LYS:HB3	1:A:193:PHE:CE2	2.54	0.43
1:B:248:GLU:HG2	3:L:19:DG:OP2	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:PHE:CZ	1:A:192:VAL:HG11	2.53	0.42
1:D:181:VAL:HG22	1:D:188:PHE:HB2	2.01	0.42
1:D:193:PHE:HE1	1:D:331:HIS:CE1	2.37	0.42
2:M:13:DG:H2'	2:M:14:DA:C8	2.54	0.42
1:A:291:MET:HB2	1:A:324:LEU:HG	2.01	0.42
1:B:153:CYS:O	1:B:154:LEU:HB2	2.19	0.42
1:C:206:ILE:CG2	1:C:230:ALA:HB2	2.49	0.42
1:B:271:GLN:HG3	1:B:285:SER:OG	2.19	0.42
1:B:311:ARG:NH1	5:B:92:HOH:O	2.52	0.42
3:N:10:DC:H2''	3:N:11:DT:O5'	2.19	0.42
1:A:253:ASN:CG	1:A:288:THR:HB	2.44	0.42
1:C:149:PHE:CD2	1:C:149:PHE:C	2.97	0.42
1:A:281:LEU:HD13	1:A:294:LEU:HG	2.01	0.42
1:A:319:LYS:O	1:A:324:LEU:HB2	2.19	0.42
1:B:151:LYS:HE3	1:B:211:ARG:CZ	2.49	0.42
1:B:154:LEU:HA	1:B:155:PRO:HD3	1.92	0.42
1:B:291:MET:HB2	1:B:324:LEU:HD12	2.01	0.42
1:C:182:ALA:HB3	1:C:238:VAL:HG21	2.02	0.42
1:C:276:LYS:HG3	1:C:277:LYS:H	1.85	0.42
3:N:15:DT:H2''	3:N:16:DG:O5'	2.19	0.42
1:C:325:GLN:HE21	1:C:327:THR:HG23	1.85	0.42
3:N:13:DT:H2''	3:N:14:DT:C6	2.55	0.42
1:D:207:ILE:CD1	1:D:209:ILE:HD11	2.50	0.42
2:M:1:DC:H2''	2:M:2:DC:H5''	2.00	0.42
2:M:4:DT:C4	2:M:5:DC:N4	2.87	0.42
1:A:201:PHE:O	1:A:202:ILE:CG1	2.68	0.42
1:C:312:LEU:CD2	1:C:334:ILE:HD12	2.49	0.42
1:B:197:LEU:HB3	1:B:201:PHE:CE2	2.54	0.42
1:B:272:LYS:CG	1:B:273:VAL:N	2.83	0.42
1:C:149:PHE:HD1	1:C:213:TYR:CD2	2.38	0.42
1:B:277:LYS:O	1:B:297:ARG:HD3	2.20	0.41
1:B:297:ARG:HE	1:B:302:MET:HE1	1.82	0.41
1:D:187:PHE:CE1	1:D:245:LYS:HE3	2.54	0.41
1:D:264:VAL:HG21	1:D:326:LEU:HD22	2.01	0.41
1:B:215:HIS:HE2	1:B:220:GLU:HB3	1.85	0.41
1:C:148:GLY:O	1:C:149:PHE:C	2.62	0.41
1:D:296:VAL:CG1	1:D:297:ARG:N	2.82	0.41
1:A:150:GLN:HB2	1:A:212:TYR:CZ	2.55	0.41
1:D:197:LEU:HD23	1:D:197:LEU:HA	1.89	0.41
1:C:251:LYS:NZ	5:C:128:HOH:O	2.50	0.41
1:D:274:THR:HB	1:D:281:LEU:HB3	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:13:DG:H2''	2:M:14:DA:H5'	2.02	0.41
1:B:298:ASN:O	1:B:302:MET:HB2	2.19	0.41
1:C:278:LYS:C	1:C:279:ASN:ND2	2.79	0.41
1:C:176:MET:CE	1:C:191:LYS:HB3	2.50	0.41
1:C:207:ILE:O	1:C:208:ILE:HD13	2.21	0.41
1:C:317:LEU:HA	1:C:326:LEU:HD23	2.03	0.41
1:A:157:MET:CE	1:A:204:LYS:HA	2.50	0.41
1:B:149:PHE:CE1	1:B:211:ARG:HB3	2.56	0.41
1:B:251:LYS:HD3	1:B:251:LYS:HA	1.72	0.41
1:C:192:VAL:HG13	1:C:221:VAL:HG11	2.03	0.41
1:C:334:ILE:HG13	1:C:335:LYS:N	2.35	0.41
1:B:184:GLU:HA	1:B:237:ASN:O	2.20	0.41
1:B:286:ASP:HB3	5:B:81:HOH:O	2.21	0.41
1:B:319:LYS:HB3	1:B:325:GLN:HB3	2.02	0.41
1:D:154:LEU:HG	1:D:156:VAL:HG13	2.03	0.41
1:D:280:ILE:C	1:D:294:LEU:HD12	2.46	0.41
1:D:286:ASP:OD1	1:D:286:ASP:C	2.63	0.41
1:A:292:GLU:O	1:A:325:GLN:HA	2.20	0.41
1:B:147:GLU:HG3	1:B:213:TYR:HE2	1.85	0.41
1:B:214:ARG:HD3	5:B:352:HOH:O	2.20	0.41
1:C:262:THR:HB	1:C:317:LEU:HB3	2.03	0.41
1:D:157:MET:HE2	1:D:159:LEU:HD23	2.03	0.41
1:D:192:VAL:HA	1:D:221:VAL:O	2.20	0.41
1:D:214:ARG:HH22	4:D:4:EDO:H21	1.86	0.41
1:D:217:GLY:HA3	4:D:6:EDO:H22	2.03	0.41
1:A:276:LYS:HB2	1:A:279:ASN:O	2.21	0.41
1:A:338:LYS:HD3	1:A:338:LYS:HA	1.61	0.41
1:B:150:GLN:HE21	1:B:150:GLN:HB3	1.63	0.41
1:B:313:THR:HG23	4:B:8:EDO:O1	2.20	0.41
1:B:338:LYS:HD3	1:B:338:LYS:HA	1.95	0.41
1:C:213:TYR:HB2	1:C:220:GLU:HB3	2.02	0.41
1:A:150:GLN:HB2	1:A:212:TYR:CE2	2.57	0.40
1:B:231:GLU:C	1:B:233:ASP:H	2.28	0.40
1:B:275:GLU:O	1:B:276:LYS:HD3	2.21	0.40
1:A:308:ASP:OD2	1:A:338:LYS:HA	2.21	0.40
1:B:293:VAL:HG23	1:B:326:LEU:C	2.46	0.40
1:D:190:VAL:HA	1:D:219:LEU:O	2.21	0.40
1:D:196:LEU:N	1:D:196:LEU:HD22	2.36	0.40
1:A:272:LYS:HG2	1:A:273:VAL:N	2.35	0.40
1:C:209:ILE:HG22	1:C:212:TYR:CD2	2.55	0.40
1:D:188:PHE:HE1	1:D:214:ARG:NH1	2.20	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:ILE:HG22	1:D:230:ALA:CB	2.51	0.40
1:A:154:LEU:HD12	1:A:155:PRO:CD	2.52	0.40
1:C:149:PHE:HA	1:C:213:TYR:C	2.46	0.40
1:C:282:PHE:HE2	1:C:295:GLY:HA3	1.86	0.40
1:D:209:ILE:HG22	1:D:212:TYR:CD2	2.57	0.40
3:L:19:DG:O3'	3:N:1:DG:H5''	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	192/208 (92%)	180 (94%)	12 (6%)	0	100	100
1	B	192/208 (92%)	176 (92%)	15 (8%)	1 (0%)	24	43
1	C	192/208 (92%)	179 (93%)	13 (7%)	0	100	100
1	D	191/208 (92%)	179 (94%)	12 (6%)	0	100	100
All	All	767/832 (92%)	714 (93%)	52 (7%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	236	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	175/185 (95%)	165 (94%)	10 (6%)	18	39
1	B	175/185 (95%)	167 (95%)	8 (5%)	24	48
1	C	175/185 (95%)	162 (93%)	13 (7%)	13	27
1	D	174/185 (94%)	168 (97%)	6 (3%)	32	60
All	All	699/740 (94%)	662 (95%)	37 (5%)	20	42

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	183	THR
1	A	209	ILE
1	A	243	ILE
1	A	254	THR
1	A	269	VAL
1	A	272	LYS
1	A	276	LYS
1	A	330	VAL
1	A	334	ILE
1	A	337	ILE
1	B	167	PHE
1	B	173	LYS
1	B	254	THR
1	B	292	GLU
1	B	299	GLU
1	B	302	MET
1	B	306	GLU
1	B	334	ILE
1	C	156	VAL
1	C	160	LYS
1	C	199	ASP
1	C	215	HIS
1	C	254	THR
1	C	260	LEU
1	C	265	ASN
1	C	293	VAL
1	C	294	LEU
1	C	317	LEU
1	C	322	GLU
1	C	327	THR
1	C	334	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	195	THR
1	D	265	ASN
1	D	271	GLN
1	D	316	THR
1	D	326	LEU
1	D	337	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	234	GLN
1	A	325	GLN
1	B	331	HIS
1	C	271	GLN
1	C	325	GLN
1	D	237	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	D	3	-	3,3,3	0.53	0	2,2,2	0.44	0
4	EDO	B	5	-	3,3,3	0.19	0	2,2,2	0.91	0
4	EDO	C	1	-	3,3,3	0.43	0	2,2,2	0.35	0
4	EDO	D	4	-	3,3,3	0.43	0	2,2,2	0.38	0
4	EDO	A	7	-	3,3,3	0.41	0	2,2,2	0.44	0
4	EDO	A	2	-	3,3,3	0.43	0	2,2,2	0.38	0
4	EDO	D	9	-	3,3,3	0.44	0	2,2,2	0.39	0
4	EDO	D	6	-	3,3,3	0.44	0	2,2,2	0.32	0
4	EDO	B	8	-	3,3,3	0.43	0	2,2,2	0.36	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	D	3	-	-	1/1/1/1	-
4	EDO	B	5	-	-	1/1/1/1	-
4	EDO	C	1	-	-	0/1/1/1	-
4	EDO	D	4	-	-	0/1/1/1	-
4	EDO	A	7	-	-	0/1/1/1	-
4	EDO	A	2	-	-	0/1/1/1	-
4	EDO	D	9	-	-	0/1/1/1	-
4	EDO	D	6	-	-	0/1/1/1	-
4	EDO	B	8	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	5	EDO	O1-C1-C2-O2
4	D	3	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	4	EDO	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	7	EDO	3	0
4	A	2	EDO	1	0
4	D	9	EDO	2	0
4	D	6	EDO	6	0
4	B	8	EDO	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	194/208 (93%)	-0.23	0 100 100	39, 69, 112, 146	0
1	B	194/208 (93%)	-0.28	1 (0%) 87 85	36, 68, 118, 148	0
1	C	194/208 (93%)	-0.28	1 (0%) 87 85	42, 69, 116, 139	0
1	D	193/208 (92%)	-0.31	1 (0%) 87 85	35, 64, 108, 131	0
2	K	19/19 (100%)	-0.34	0 100 100	45, 65, 153, 175	0
2	M	19/19 (100%)	-0.06	0 100 100	63, 84, 125, 126	0
3	L	19/19 (100%)	0.12	1 (5%) 32 28	57, 102, 188, 192	0
3	N	19/19 (100%)	-0.19	0 100 100	61, 77, 118, 118	0
All	All	851/908 (93%)	-0.26	4 (0%) 87 85	35, 69, 120, 192	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	212	TYR	2.5
3	L	1	DG	2.4
1	D	148	GLY	2.3
1	B	237	ASN	2.1

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	A	2	4/4	0.95	0.12	53,54,55,55	0
4	EDO	D	4	4/4	0.96	0.08	49,52,60,64	0
4	EDO	D	9	4/4	0.96	0.20	54,57,62,68	0
4	EDO	D	3	4/4	0.97	0.06	39,40,42,45	0
4	EDO	B	5	4/4	0.97	0.07	52,53,57,60	0
4	EDO	D	6	4/4	0.97	0.07	52,60,64,68	0
4	EDO	B	8	4/4	0.97	0.09	44,44,50,54	0
4	EDO	A	7	4/4	0.98	0.13	39,42,43,49	0
4	EDO	C	1	4/4	0.99	0.08	39,44,47,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.