



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:42 AM UTC

PDB ID : 8AMV / pdb_00008amv
Title : RepB pMV158 hexamer
Authors : Machon, C.; Amodio, J.; Boer, R.D.; Ruiz-Maso, J.A.; del Solar, G.; Coll, M.
Deposited on : 2022-08-04
Resolution : 2.77 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

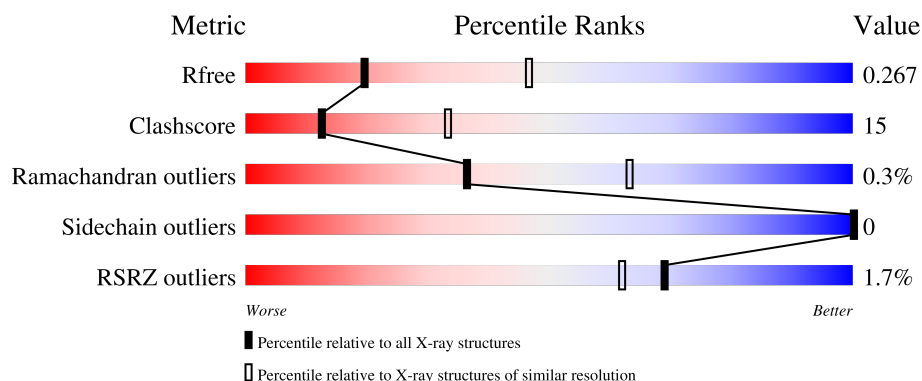
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	210	 74% 21% 5%
1	B	210	 63% 30% 5%
1	C	210	 61% 32% 5%
1	D	210	 7% 42% 37% 20%
1	E	210	 73% 21% 5%

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Mol	Chain	Length	Quality of chain
1	F	210	% 77%18%5%

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
2	PO4	B	302	-	-	X	-

2 Entry composition [i](#)

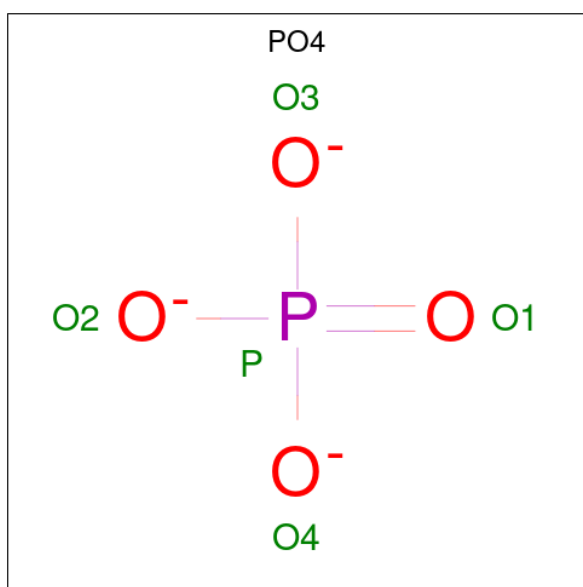
There are 4 unique types of molecules in this entry. The entry contains 9615 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Replication protein RepB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1631	1060	266	301	4			
1	B	200	Total	C	N	O	S	0	0	0
			1631	1060	266	301	4			
1	C	201	Total	C	N	O	S	0	0	0
			1640	1066	268	302	4			
1	D	167	Total	C	N	O	S	0	0	0
			1360	888	220	248	4			
1	E	201	Total	C	N	O	S	0	0	0
			1640	1066	268	302	4			
1	F	200	Total	C	N	O	S	0	0	0
			1631	1060	266	301	4			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Na 1 1	0	0

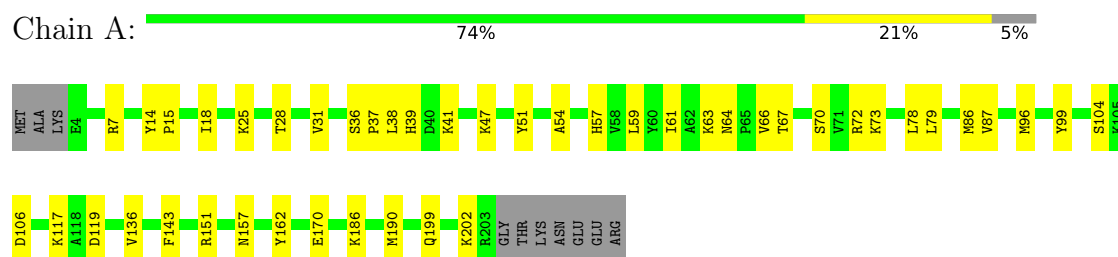
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	20	Total O 20 20	0	0
4	B	8	Total O 8 8	0	0
4	D	3	Total O 3 3	0	0
4	E	6	Total O 6 6	0	0
4	F	14	Total O 14 14	0	0

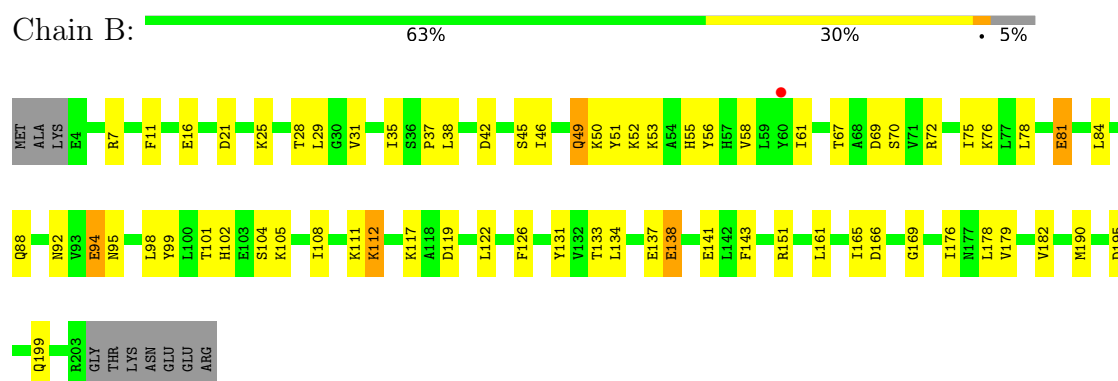
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

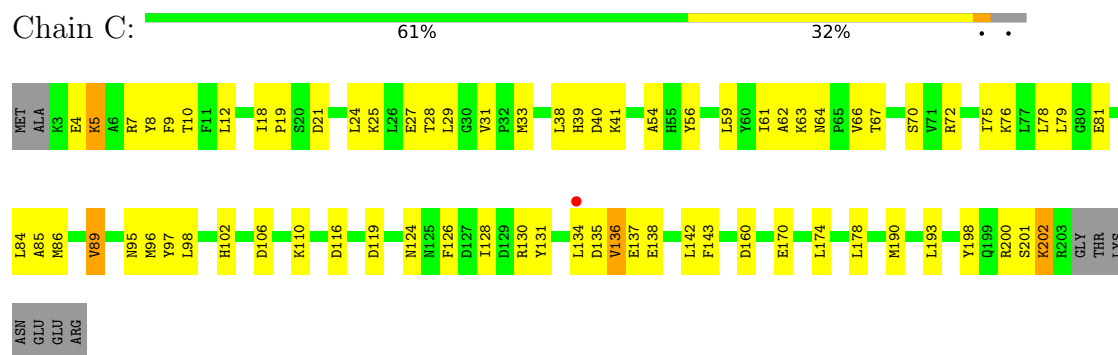
• Molecule 1: Replication protein RepB



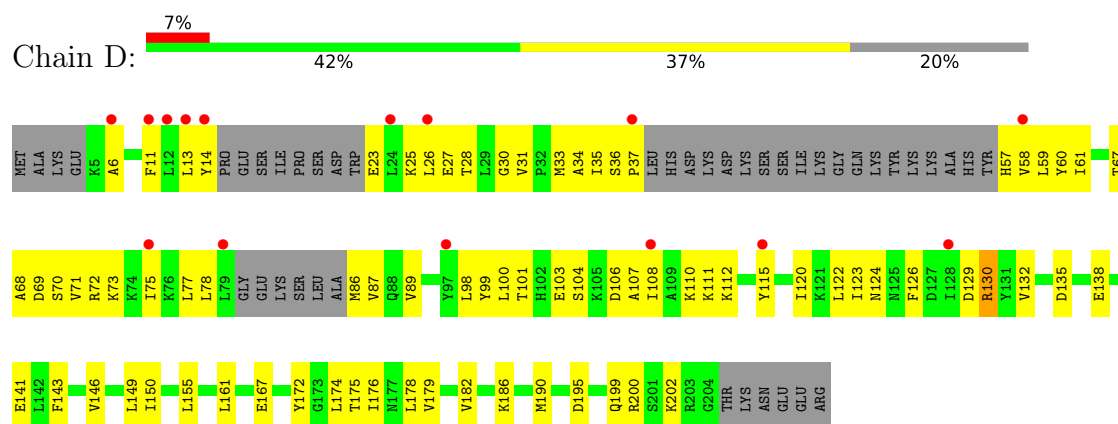
• Molecule 1: Replication protein RepB



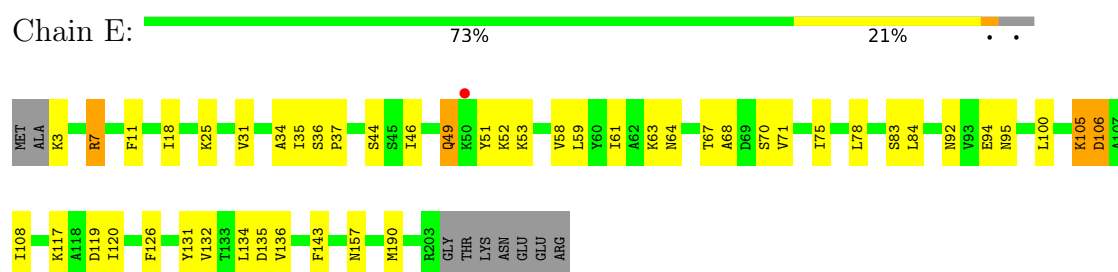
• Molecule 1: Replication protein RepB



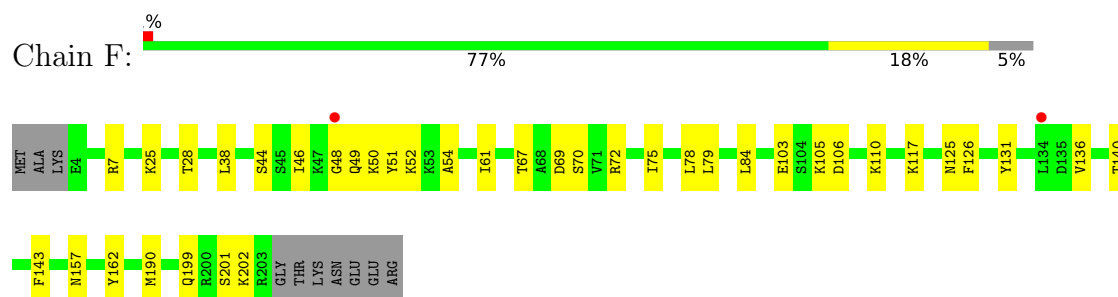
• Molecule 1: Replication protein RepB



• Molecule 1: Replication protein RepB



• Molecule 1: Replication protein RepB



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.91Å 79.33Å 114.90Å 90.00° 104.92° 90.00°	Depositor
Resolution (Å)	45.48 – 2.77 45.48 – 2.77	Depositor EDS
% Data completeness (in resolution range)	99.2 (45.48-2.77) 99.2 (45.48-2.77)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.77Å)	Xtriage
Refinement program	REFMAC 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.218 , 0.267 0.221 , 0.267	Depositor DCC
R_{free} test set	1976 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	76.7	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9615	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, PO4

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	0.50	0/1663	0.81	1/2244 (0.0%)
1	B	0.49	0/1663	0.87	4/2244 (0.2%)
1	C	0.71	2/1672 (0.1%)	1.01	14/2255 (0.6%)
1	D	0.50	1/1381 (0.1%)	0.75	0/1861
1	E	0.45	0/1672	0.84	8/2255 (0.4%)
1	F	0.55	0/1663	0.82	0/2244
All	All	0.54	3/9714 (0.0%)	0.86	27/13103 (0.2%)

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	B	0	1
1	E	0	2
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	18	ILE	C-O	-19.30	1.10	1.24
1	C	18	ILE	CA-C	8.38	1.61	1.52
1	D	30	GLY	C-N	5.30	1.39	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	18	ILE	O-C-N	13.89	130.44	121.69
1	C	202	LYS	CB-CA-C	-9.70	96.03	111.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	94	GLU	CA-CB-CG	7.37	128.83	114.10
1	E	106	ASP	N-CA-CB	-6.75	99.82	110.28
1	C	18	ILE	CA-C-N	-6.35	113.12	119.92
1	C	18	ILE	C-N-CA	-6.35	113.12	119.92
1	E	49	GLN	CB-CA-C	6.06	119.20	109.02
1	B	49	GLN	CA-CB-CG	5.97	126.05	114.10
1	E	49	GLN	CA-CB-CG	5.96	126.03	114.10
1	C	61	ILE	CA-CB-CG1	5.94	120.49	110.40
1	C	201	SER	CA-C-N	-5.90	113.75	122.83
1	C	201	SER	C-N-CA	-5.90	113.75	122.83
1	E	7	ARG	CD-NE-CZ	5.76	132.47	124.40
1	C	18	ILE	N-CA-C	-5.71	104.50	109.19
1	E	105	LYS	CA-C-N	5.47	128.63	120.31
1	E	105	LYS	C-N-CA	5.47	128.63	120.31
1	C	5	LYS	CD-CE-NZ	-5.46	94.42	111.90
1	C	135	ASP	CA-C-N	5.29	131.49	121.97
1	C	135	ASP	C-N-CA	5.29	131.49	121.97
1	B	112	LYS	CD-CE-NZ	-5.29	94.99	111.90
1	B	81	GLU	CA-CB-CG	-5.25	103.59	114.10
1	A	117	LYS	N-CA-CB	-5.21	101.73	110.39
1	C	202	LYS	CD-CE-NZ	5.21	128.57	111.90
1	E	3	LYS	CA-C-N	5.15	128.16	120.95
1	E	3	LYS	C-N-CA	5.15	128.16	120.95
1	C	89	VAL	CA-C-N	-5.10	115.61	122.91
1	C	89	VAL	C-N-CA	-5.10	115.61	122.91

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	138	GLU	Sidechain
1	E	106	ASP	Sidechain
1	E	7	ARG	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	A	1631	0	1673	36	1
1	B	1631	0	1673	71	1
1	C	1640	0	1686	73	0
1	D	1360	0	1403	70	0
1	E	1640	0	1685	31	0
1	F	1631	0	1673	28	0
2	A	5	0	0	0	0
2	B	5	0	0	6	0
2	C	5	0	0	1	0
2	D	5	0	0	1	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	B	1	0	0	0	0
4	A	20	0	0	0	0
4	B	8	0	0	0	0
4	D	3	0	0	0	0
4	E	6	0	0	0	0
4	F	14	0	0	1	0
All	All	9615	0	9793	297	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (297) 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:B:302:PO4:O4	1:C:130:ARG:NH1	1.75	1.18
1:B:67:THR:HG23	1:B:70:SER:HB3	1.25	1.14
1:E:46:ILE:HD11	1:E:49:GLN:OE1	1.53	1.05
1:E:31:VAL:HG13	1:E:61:ILE:O	1.61	1.00
1:C:134:LEU:HG	1:C:138:GLU:HG2	1.50	0.93
1:D:108:ILE:HA	1:D:111:LYS:HG2	1.53	0.90
1:C:38:LEU:HD21	1:C:54:ALA:HB1	1.54	0.89
1:A:67:THR:HG23	1:A:70:SER:HB3	1.54	0.89
1:F:67:THR:HG23	1:F:70:SER:H	1.38	0.89
1:D:13:LEU:HD23	1:D:14:TYR:H	1.39	0.88
1:C:198:TYR:CE2	1:C:202:LYS:NZ	2.41	0.88
1:C:84:LEU:HD13	1:C:85:ALA:N	1.89	0.88
1:F:49:GLN:OE1	1:F:51:TYR:O	1.92	0.88
1:B:81:GLU:O	1:B:81:GLU:HG3	1.75	0.86
1:D:13:LEU:HD23	1:D:14:TYR:N	1.90	0.86
1:C:136:VAL:HG13	1:C:137:GLU:OE1	1.75	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:VAL:HG12	1:C:70:SER:OG	1.77	0.84
1:C:75:ILE:O	1:C:79:LEU:HD12	1.79	0.82
1:B:72:ARG:HG2	1:B:76:LYS:HD2	1.61	0.82
2:B:302:PO4:O4	1:C:130:ARG:CZ	2.30	0.79
1:D:106:ASP:HB3	1:E:134:LEU:HD11	1.64	0.78
1:D:104:SER:H	1:D:107:ALA:HB3	1.49	0.77
1:C:5:LYS:HG2	1:C:67:THR:HG22	1.69	0.75
1:F:38:LEU:HD11	1:F:54:ALA:HB1	1.69	0.75
1:D:178:LEU:O	1:D:182:VAL:HG23	1.87	0.75
1:B:46:ILE:CD1	1:B:49:GLN:OE1	2.35	0.74
1:B:67:THR:CG2	1:B:70:SER:HB3	2.11	0.73
1:D:122:LEU:HD23	1:D:126:PHE:HD2	1.54	0.73
1:F:25:LYS:HB3	1:F:78:LEU:HD11	1.69	0.73
1:B:61:ILE:HD13	1:B:131:TYR:CZ	2.23	0.72
1:D:26:LEU:HD23	1:D:33:MET:HE1	1.70	0.72
1:B:178:LEU:O	1:B:182:VAL:HG23	1.90	0.71
1:B:134:LEU:O	1:B:138:GLU:HG3	1.91	0.70
1:B:46:ILE:HD11	1:B:49:GLN:HB3	1.73	0.70
1:C:75:ILE:HG13	1:C:79:LEU:HD11	1.73	0.70
1:A:14:TYR:O	1:A:18:ILE:HG12	1.92	0.70
1:B:67:THR:HG23	1:B:70:SER:CB	2.14	0.70
1:B:16:GLU:OE2	1:B:53:LYS:NZ	2.22	0.70
1:E:18:ILE:HG22	1:E:83:SER:HB2	1.74	0.69
1:F:103:GLU:OE2	1:F:117:LYS:HE2	1.91	0.69
1:B:61:ILE:CD1	1:B:131:TYR:CZ	2.77	0.68
1:B:102:HIS:NE2	1:B:112:LYS:HE3	2.09	0.68
1:D:57:HIS:NE2	2:D:301:PO4:O3	2.27	0.67
1:F:44:SER:OG	1:F:46:ILE:HG22	1.94	0.67
1:C:106:ASP:OD1	1:C:110:LYS:HE2	1.96	0.66
1:A:96:MET:HE3	1:A:99:TYR:HB3	1.78	0.65
1:B:61:ILE:HD13	1:B:131:TYR:CE2	2.31	0.65
1:D:6:ALA:HB3	1:D:89:VAL:HG23	1.78	0.65
1:C:75:ILE:HG13	1:C:79:LEU:CD1	2.26	0.65
1:C:29:LEU:HD21	1:C:78:LEU:HG	1.77	0.64
1:B:133:THR:O	1:B:137:GLU:HG2	1.97	0.64
1:D:68:ALA:HA	1:D:71:VAL:HG12	1.79	0.64
2:B:302:PO4:P	1:C:130:ARG:HH22	2.21	0.64
1:D:26:LEU:H	1:D:26:LEU:HD12	1.63	0.63
1:B:72:ARG:O	1:B:76:LYS:HG3	1.99	0.63
1:D:11:PHE:HB3	1:D:87:VAL:HG22	1.79	0.63
1:A:25:LYS:HB3	1:A:78:LEU:HD11	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:PRO:HB3	1:A:119:ASP:HB2	1.81	0.63
1:C:75:ILE:HD12	1:C:78:LEU:HB3	1.79	0.62
1:A:72:ARG:HH12	1:A:73:LYS:HE3	1.65	0.62
1:C:84:LEU:HD13	1:C:85:ALA:C	2.25	0.62
1:E:157:ASN:HD22	1:F:199:GLN:NE2	1.98	0.62
1:F:49:GLN:O	1:F:51:TYR:N	2.29	0.61
1:D:67:THR:HG23	1:D:70:SER:H	1.65	0.61
1:B:7:ARG:NH2	1:B:138:GLU:OE1	2.32	0.61
1:B:42:ASP:HB3	1:B:52:LYS:HD2	1.81	0.61
1:B:46:ILE:HD11	1:B:49:GLN:OE1	2.00	0.60
1:D:34:ALA:HB2	1:D:126:PHE:CE2	2.37	0.60
1:B:111:LYS:HD3	1:B:112:LYS:N	2.17	0.60
1:D:69:ASP:O	1:D:72:ARG:HG3	2.02	0.60
1:C:75:ILE:CG1	1:C:79:LEU:HD11	2.32	0.60
1:C:84:LEU:HD22	1:C:85:ALA:H	1.66	0.60
1:F:143:PHE:HE1	1:F:190:MET:HE3	1.66	0.60
1:C:66:VAL:CG1	1:C:70:SER:OG	2.48	0.59
1:F:143:PHE:CE1	1:F:190:MET:HE3	2.37	0.59
1:E:31:VAL:CG1	1:E:61:ILE:O	2.44	0.59
1:A:64:ASN:O	1:A:66:VAL:HG23	2.02	0.59
1:B:49:GLN:O	1:B:51:TYR:N	2.35	0.59
1:D:122:LEU:HD23	1:D:126:PHE:CD2	2.38	0.59
1:F:105:LYS:HD2	1:F:105:LYS:N	2.18	0.59
1:D:31:VAL:HG11	1:D:60:TYR:CZ	2.38	0.59
1:F:106:ASP:OD1	1:F:110:LYS:NZ	2.36	0.59
1:C:75:ILE:CD1	1:C:79:LEU:HD11	2.33	0.58
1:C:19:PRO:HB2	1:C:21:ASP:OD1	2.04	0.58
1:D:13:LEU:CD2	1:D:14:TYR:N	2.64	0.58
1:D:101:THR:OG1	1:D:103:GLU:HB2	2.02	0.58
1:B:72:ARG:HH11	1:B:76:LYS:NZ	2.02	0.57
1:C:84:LEU:HD13	1:C:85:ALA:H	1.68	0.57
1:B:141:GLU:O	1:B:141:GLU:HG3	2.03	0.57
1:F:44:SER:HB2	1:F:52:LYS:HG3	1.87	0.57
1:C:38:LEU:HD12	1:C:56:TYR:CZ	2.39	0.56
1:D:25:LYS:O	1:D:78:LEU:HD11	2.05	0.56
1:E:35:ILE:HD12	1:E:58:VAL:HG22	1.87	0.56
1:B:102:HIS:O	1:B:108:ILE:HD13	2.05	0.56
1:C:25:LYS:O	1:C:28:THR:HG22	2.04	0.56
1:C:4:GLU:HG3	1:C:170:GLU:HG3	1.86	0.56
1:E:135:ASP:OD1	1:E:136:VAL:N	2.39	0.56
1:A:170:GLU:H	1:A:170:GLU:CD	2.14	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:25:LYS:O	1:F:28:THR:HB	2.05	0.55
1:A:72:ARG:NH1	1:A:73:LYS:HE3	2.22	0.55
1:E:44:SER:OG	1:E:46:ILE:HG12	2.05	0.55
1:A:86:MET:HE3	1:A:87:VAL:H	1.71	0.55
1:E:44:SER:HB2	1:E:52:LYS:HG3	1.88	0.55
1:E:94:GLU:HG3	1:E:132:VAL:HG12	1.89	0.54
1:B:61:ILE:O	1:B:61:ILE:HG13	2.06	0.54
1:C:25:LYS:HG2	1:C:78:LEU:HD21	1.89	0.54
1:C:84:LEU:HD11	1:C:86:MET:O	2.07	0.54
1:C:29:LEU:HB3	1:C:31:VAL:HG23	1.90	0.54
1:D:31:VAL:HG11	1:D:60:TYR:CE2	2.42	0.54
1:D:67:THR:CG2	1:D:70:SER:HB3	2.37	0.54
1:E:61:ILE:HD13	1:E:131:TYR:CZ	2.43	0.53
1:C:9:PHE:CE1	1:C:89:VAL:HG22	2.44	0.53
1:B:38:LEU:HD13	1:B:56:TYR:CZ	2.44	0.53
1:D:132:VAL:HA	1:D:135:ASP:OD1	2.09	0.52
1:D:13:LEU:HD22	1:D:14:TYR:O	2.08	0.52
1:A:25:LYS:HB3	1:A:78:LEU:CD1	2.40	0.52
1:A:7:ARG:NH2	1:A:63:LYS:O	2.42	0.52
1:B:46:ILE:HD12	1:B:49:GLN:OE1	2.07	0.52
1:D:146:VAL:O	1:D:150:ILE:HG13	2.10	0.52
1:A:15:PRO:HD2	1:A:54:ALA:O	2.09	0.52
1:A:25:LYS:O	1:A:28:THR:HB	2.10	0.52
1:D:23:GLU:HA	1:D:26:LEU:HD13	1.92	0.52
1:D:195:ASP:O	1:D:199:GLN:HG2	2.10	0.52
1:D:138:GLU:HA	1:D:141:GLU:HB2	1.92	0.51
1:F:201:SER:C	1:F:202:LYS:HD3	2.35	0.51
1:C:143:PHE:CE1	1:C:190:MET:HE3	2.45	0.51
1:A:31:VAL:HG13	1:A:61:ILE:O	2.10	0.51
1:D:35:ILE:HG12	1:D:58:VAL:HG22	1.93	0.51
1:D:13:LEU:CD2	1:D:14:TYR:O	2.59	0.51
1:B:25:LYS:O	1:B:78:LEU:HD11	2.10	0.51
1:A:151:ARG:HD2	1:F:162:TYR:CD2	2.45	0.51
1:C:33:MET:HA	1:C:59:LEU:O	2.11	0.51
1:F:75:ILE:HG23	1:F:79:LEU:HD12	1.93	0.51
1:A:202:LYS:HD3	1:A:202:LYS:O	2.10	0.50
1:B:29:LEU:HD21	1:B:78:LEU:HG	1.94	0.50
1:B:76:LYS:HG2	1:B:84:LEU:HD21	1.93	0.50
2:B:302:PO4:P	1:C:130:ARG:HH12	2.30	0.50
1:C:28:THR:HG22	1:C:29:LEU:HD22	1.94	0.50
1:C:142:LEU:HD22	1:C:174:LEU:HD21	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:ASN:ND2	1:B:104:SER:HB2	2.27	0.50
1:B:143:PHE:HE1	1:B:190:MET:HE3	1.76	0.50
1:C:38:LEU:HD12	1:C:56:TYR:CE2	2.47	0.50
1:A:143:PHE:CE1	1:A:190:MET:HE3	2.47	0.50
1:B:61:ILE:CD1	1:B:131:TYR:CE1	2.94	0.50
1:B:101:THR:HG21	1:B:117:LYS:HG2	1.94	0.50
2:B:302:PO4:O4	1:C:130:ARG:NH2	2.45	0.50
1:F:61:ILE:HD13	1:F:131:TYR:CZ	2.47	0.50
1:E:126:PHE:CD1	1:E:126:PHE:C	2.90	0.49
1:B:88:GLN:NE2	1:C:137:GLU:HG3	2.27	0.49
1:A:104:SER:OG	1:A:106:ASP:OD1	2.28	0.49
1:D:37:PRO:HG2	1:D:115:TYR:HB3	1.94	0.49
1:B:21:ASP:OD1	1:B:25:LYS:HE2	2.12	0.49
1:D:175:THR:O	1:D:179:VAL:HG13	2.13	0.49
1:C:116:ASP:HB3	1:C:119:ASP:OD1	2.13	0.49
1:A:199:GLN:NE2	1:F:157:ASN:HD22	2.10	0.48
1:B:7:ARG:CZ	1:B:138:GLU:OE1	2.61	0.48
1:B:165:ILE:O	1:B:169:GLY:N	2.46	0.48
1:D:72:ARG:NH2	1:D:86:MET:HA	2.29	0.48
1:D:6:ALA:HB3	1:D:89:VAL:CG2	2.44	0.48
1:B:55:HIS:NE2	2:B:302:PO4:O1	2.42	0.48
1:A:162:TYR:CD2	1:B:151:ARG:HD2	2.49	0.48
1:C:160:ASP:OD1	1:D:200:ARG:NH2	2.46	0.48
1:D:69:ASP:O	1:D:73:LYS:HG2	2.13	0.48
1:D:31:VAL:N	1:D:124:ASN:OD1	2.47	0.47
1:C:9:PHE:HZ	1:C:66:VAL:O	1.97	0.47
1:D:143:PHE:CE1	1:D:190:MET:HE3	2.49	0.47
1:C:75:ILE:HD11	1:C:79:LEU:HD11	1.95	0.47
1:E:63:LYS:HG3	1:E:64:ASN:N	2.30	0.47
1:B:161:LEU:O	1:B:165:ILE:HG12	2.14	0.47
1:B:11:PHE:CZ	1:B:75:ILE:HD13	2.50	0.47
1:C:174:LEU:HD23	1:C:178:LEU:HD23	1.96	0.47
1:C:97:TYR:CG	1:C:128:ILE:HD12	2.50	0.47
1:F:125:ASN:ND2	4:F:401:HOH:O	2.47	0.47
1:A:41:LYS:HG2	1:A:51:TYR:CE2	2.49	0.46
1:B:35:ILE:HG12	1:B:58:VAL:HG22	1.96	0.46
1:C:59:LEU:HD21	1:C:96:MET:HG2	1.96	0.46
1:C:102:HIS:NE2	2:C:301:PO4:O2	2.35	0.46
1:F:69:ASP:N	1:F:69:ASP:OD1	2.46	0.46
1:B:94:GLU:O	1:B:98:LEU:HD12	2.15	0.46
1:A:38:LEU:HD11	1:A:54:ALA:HB1	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:LYS:H	1:A:47:LYS:HD3	1.80	0.46
1:A:157:ASN:HB3	1:B:199:GLN:OE1	2.16	0.46
1:E:117:LYS:HG2	1:E:120:ILE:HD12	1.97	0.46
1:C:134:LEU:CG	1:C:138:GLU:HG2	2.34	0.46
1:E:143:PHE:CE1	1:E:190:MET:HE3	2.50	0.46
1:A:47:LYS:HE2	1:A:47:LYS:HB2	1.51	0.46
1:E:37:PRO:HB3	1:E:119:ASP:HB2	1.98	0.46
1:B:195:ASP:O	1:B:199:GLN:HG3	2.16	0.46
1:D:67:THR:HG22	1:D:70:SER:HB3	1.97	0.46
1:D:150:ILE:HG23	1:D:155:LEU:HB2	1.97	0.45
1:D:98:LEU:HB3	1:D:103:GLU:O	2.16	0.45
1:E:36:SER:HB3	1:E:100:LEU:HD22	1.98	0.45
1:E:49:GLN:NE2	1:E:51:TYR:O	2.48	0.45
1:E:67:THR:O	1:E:70:SER:HB3	2.15	0.45
1:B:61:ILE:HD11	1:B:131:TYR:CZ	2.50	0.45
1:C:31:VAL:N	1:C:124:ASN:HD22	2.15	0.45
1:D:36:SER:OG	1:D:100:LEU:HB3	2.17	0.45
1:D:59:LEU:HD23	1:D:60:TYR:N	2.31	0.45
1:C:190:MET:HE2	1:C:193:LEU:HD12	1.98	0.45
1:B:46:ILE:CD1	1:B:49:GLN:HB3	2.44	0.45
1:B:72:ARG:CG	1:B:76:LYS:HD2	2.41	0.45
1:A:36:SER:OG	1:A:37:PRO:O	2.35	0.45
1:D:174:LEU:HD22	1:D:179:VAL:HG12	1.98	0.45
1:F:72:ARG:HG3	1:F:84:LEU:HD11	1.99	0.45
1:A:67:THR:CG2	1:A:70:SER:HB3	2.34	0.45
1:C:160:ASP:OD1	1:D:200:ARG:NH1	2.49	0.45
1:D:25:LYS:O	1:D:28:THR:HB	2.16	0.45
1:E:53:LYS:HB3	1:E:53:LYS:HZ3	1.81	0.45
1:B:108:ILE:HD12	1:B:108:ILE:HA	1.89	0.45
1:C:10:THR:HG21	1:C:96:MET:HE3	1.99	0.45
1:C:39:HIS:O	1:C:54:ALA:HA	2.16	0.45
1:D:176:ILE:O	1:D:179:VAL:HG22	2.17	0.45
1:A:59:LEU:HD22	1:A:61:ILE:HD11	1.98	0.44
1:A:143:PHE:HE1	1:A:190:MET:HE3	1.82	0.44
1:B:72:ARG:NH1	1:B:76:LYS:NZ	2.66	0.44
1:C:29:LEU:HB3	1:C:31:VAL:CG2	2.47	0.44
1:C:72:ARG:O	1:C:76:LYS:HG3	2.17	0.44
1:A:136:VAL:HG22	1:F:7:ARG:NH1	2.32	0.44
1:B:99:TYR:CE1	1:B:112:LYS:HE2	2.52	0.44
1:D:108:ILE:HA	1:D:111:LYS:CG	2.36	0.44
1:B:45:SER:O	1:B:45:SER:OG	2.34	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:ASP:O	1:B:72:ARG:HB3	2.17	0.44
1:C:40:ASP:OD1	1:C:41:LYS:HG2	2.18	0.44
1:C:63:LYS:HD2	1:C:63:LYS:HA	1.79	0.44
1:D:202:LYS:HD2	1:D:202:LYS:HA	1.67	0.44
1:B:143:PHE:CE1	1:B:190:MET:HE3	2.52	0.44
1:C:24:LEU:HA	1:C:27:GLU:OE1	2.18	0.44
1:B:72:ARG:HH11	1:B:76:LYS:HZ2	1.64	0.44
1:F:44:SER:N	1:F:50:LYS:O	2.35	0.44
1:C:75:ILE:HD12	1:C:75:ILE:HA	1.79	0.44
1:B:21:ASP:CG	1:B:25:LYS:HE2	2.43	0.43
1:D:73:LYS:O	1:D:77:LEU:HD23	2.18	0.43
1:A:78:LEU:C	1:A:78:LEU:HD13	2.43	0.43
1:B:37:PRO:HB3	1:B:119:ASP:CB	2.48	0.43
1:D:167:GLU:HA	1:D:167:GLU:OE1	2.18	0.43
1:F:126:PHE:CD1	1:F:126:PHE:C	2.96	0.43
1:A:37:PRO:O	1:A:39:HIS:HD2	2.02	0.43
1:B:95:ASN:HD22	1:B:104:SER:HB2	1.83	0.43
1:D:155:LEU:HD13	1:D:161:LEU:HA	1.99	0.43
1:E:59:LEU:HD23	1:E:59:LEU:HA	1.84	0.43
1:F:136:VAL:O	1:F:140:THR:HG23	2.18	0.43
1:C:170:GLU:H	1:C:170:GLU:CD	2.25	0.43
1:D:99:TYR:C	1:D:101:THR:N	2.76	0.43
1:B:53:LYS:HE3	1:C:126:PHE:O	2.19	0.43
1:C:12:LEU:O	1:C:84:LEU:HD23	2.19	0.43
1:B:76:LYS:HG2	1:B:84:LEU:CD2	2.48	0.43
1:B:92:ASN:HD21	1:B:94:GLU:HB2	1.83	0.43
1:D:34:ALA:HB2	1:D:126:PHE:CD2	2.54	0.43
1:E:34:ALA:HB3	1:E:59:LEU:HB2	2.00	0.43
1:B:25:LYS:C	1:B:78:LEU:HD11	2.44	0.43
1:C:95:ASN:HA	1:C:98:LEU:HD12	2.01	0.43
1:E:92:ASN:HD22	1:E:95:ASN:ND2	2.17	0.43
1:C:200:ARG:HD3	1:C:200:ARG:HA	1.54	0.42
1:E:84:LEU:H	1:E:84:LEU:HD23	1.82	0.42
1:C:7:ARG:HG2	1:C:62:ALA:O	2.18	0.42
1:C:76:LYS:HD3	1:C:81:GLU:OE2	2.19	0.42
1:D:129:ASP:O	1:D:130:ARG:C	2.62	0.42
1:B:37:PRO:HB3	1:B:119:ASP:HB3	2.02	0.42
1:C:137:GLU:CD	1:C:137:GLU:H	2.26	0.42
1:D:31:VAL:HG22	1:D:61:ILE:O	2.19	0.42
1:D:33:MET:SD	1:D:123:ILE:HD11	2.60	0.42
1:F:61:ILE:HD11	1:F:126:PHE:CE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:LEU:HD22	1:B:126:PHE:CD1	2.55	0.42
1:D:75:ILE:O	1:D:78:LEU:HB3	2.20	0.42
1:C:8:TYR:CE1	1:C:131:TYR:HB3	2.55	0.42
1:C:64:ASN:O	1:C:66:VAL:HG23	2.19	0.42
1:C:128:ILE:HA	1:C:131:TYR:CD2	2.54	0.42
1:A:36:SER:HB3	1:A:57:HIS:H	1.84	0.42
1:F:67:THR:O	1:F:70:SER:HB3	2.19	0.42
1:B:88:GLN:HE21	1:C:137:GLU:HG3	1.85	0.41
1:B:102:HIS:CD2	1:B:112:LYS:HE3	2.55	0.41
1:D:110:LYS:HE3	1:D:112:LYS:HG3	2.02	0.41
1:B:25:LYS:O	1:B:28:THR:HB	2.20	0.41
1:D:36:SER:HB2	1:D:57:HIS:O	2.20	0.41
1:C:67:THR:O	1:C:70:SER:OG	2.38	0.41
1:D:100:LEU:N	1:D:100:LEU:HD23	2.35	0.41
1:C:75:ILE:CD1	1:C:79:LEU:CD1	2.98	0.41
1:D:26:LEU:HD22	1:D:35:ILE:HD11	2.01	0.41
1:D:27:GLU:HG2	1:D:123:ILE:HD12	2.02	0.41
1:E:105:LYS:HA	1:E:108:ILE:CG1	2.51	0.41
1:B:31:VAL:HG13	1:B:61:ILE:O	2.21	0.41
1:B:176:ILE:O	1:B:179:VAL:HG22	2.21	0.41
1:C:5:LYS:CG	1:C:67:THR:HG22	2.47	0.41
1:D:143:PHE:CE1	1:D:186:LYS:HD3	2.55	0.41
1:A:78:LEU:HD12	1:A:79:LEU:HG	2.02	0.41
1:A:143:PHE:CE1	1:A:186:LYS:HD2	2.56	0.41
1:B:105:LYS:O	1:B:108:ILE:HG22	2.21	0.41
1:D:25:LYS:HB3	1:D:78:LEU:HD21	2.03	0.41
1:E:11:PHE:CZ	1:E:75:ILE:HD13	2.55	0.41
1:E:143:PHE:HE1	1:E:190:MET:HE3	1.85	0.41
1:D:35:ILE:O	1:D:120:ILE:HG23	2.20	0.40
1:D:149:LEU:HD21	1:D:172:TYR:CE2	2.57	0.40
1:E:68:ALA:O	1:E:71:VAL:HG22	2.20	0.40
1:D:107:ALA:C	1:D:111:LYS:HD2	2.46	0.40
1:E:134:LEU:HD12	1:E:135:ASP:H	1.85	0.40
1:D:70:SER:OG	1:D:71:VAL:N	2.54	0.40
1:E:25:LYS:HB3	1:E:78:LEU:HD11	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:LYS:NZ	1:B:166:ASP:OD2[2_545]	2.16	0.04

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	198/210 (94%)	195 (98%)	3 (2%)	0	100	100
1	B	198/210 (94%)	190 (96%)	7 (4%)	1 (0%)	24	52
1	C	199/210 (95%)	193 (97%)	5 (2%)	1 (0%)	24	52
1	D	159/210 (76%)	153 (96%)	5 (3%)	1 (1%)	21	47
1	E	199/210 (95%)	198 (100%)	1 (0%)	0	100	100
1	F	198/210 (94%)	193 (98%)	4 (2%)	1 (0%)	24	52
All	All	1151/1260 (91%)	1122 (98%)	25 (2%)	4 (0%)	36	63

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	136	VAL
1	B	50	LYS
1	F	48	GLY
1	D	130	ARG

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	180/188 (96%)	180 (100%)	0	100	100
1	B	180/188 (96%)	180 (100%)	0	100	100
1	C	181/188 (96%)	181 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	150/188 (80%)	150 (100%)	0	100	100
1	E	181/188 (96%)	181 (100%)	0	100	100
1	F	180/188 (96%)	180 (100%)	0	100	100
All	All	1052/1128 (93%)	1052 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	88	GLN
1	A	113	HIS
1	A	199	GLN
1	B	39	HIS
1	B	88	GLN
1	B	92	ASN
1	B	95	ASN
1	B	124	ASN
1	C	156	GLN
1	D	156	GLN
1	D	168	ASN
1	E	92	ASN
1	E	180	ASN
1	F	95	ASN
1	F	168	ASN
1	F	180	ASN
1	F	199	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	D	301	-	4,4,4	1.04	0	6,6,6	0.77	0
2	PO4	F	301	-	4,4,4	1.18	0	6,6,6	1.03	0
2	PO4	A	301	-	4,4,4	1.21	0	6,6,6	0.82	0
2	PO4	C	301	-	4,4,4	1.00	0	6,6,6	0.63	0
2	PO4	E	301	-	4,4,4	0.92	0	6,6,6	0.62	0
2	PO4	B	302	-	4,4,4	0.85	0	6,6,6	0.61	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.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	PO4	1	0
2	C	301	PO4	1	0
2	B	302	PO4	6	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	200/210 (95%)	-0.25	0 100 100	41, 70, 98, 113	0
1	B	200/210 (95%)	0.15	1 (0%) 87 83	49, 88, 134, 159	0
1	C	201/210 (95%)	0.19	1 (0%) 87 83	56, 101, 138, 169	0
1	D	167/210 (79%)	0.59	15 (8%) 15 12	57, 134, 167, 195	0
1	E	201/210 (95%)	-0.19	1 (0%) 87 83	50, 78, 113, 121	0
1	F	200/210 (95%)	-0.35	2 (1%) 79 74	40, 57, 104, 137	0
All	All	1169/1260 (92%)	0.01	20 (1%) 69 62	40, 82, 145, 195	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	134	LEU	3.4
1	D	79	LEU	3.2
1	D	58	VAL	3.0
1	D	75	ILE	2.9
1	D	115	TYR	2.9
1	D	26	LEU	2.9
1	D	11	PHE	2.8
1	D	108	ILE	2.7
1	D	13	LEU	2.6
1	D	97	TYR	2.6
1	E	50	LYS	2.5
1	D	24	LEU	2.3
1	C	134	LEU	2.3
1	B	60	TYR	2.2
1	D	37	PRO	2.2
1	D	14	TYR	2.1
1	D	128	ILE	2.1
1	F	48	GLY	2.0
1	D	6	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	12	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
3	NA	B	301	1/1	0.24	0.29	148,148,148,148	0
2	PO4	D	301	5/5	0.77	0.07	150,154,182,192	0
2	PO4	C	301	5/5	0.93	0.07	87,88,97,100	0
2	PO4	B	302	5/5	0.96	0.06	87,89,97,102	0
2	PO4	E	301	5/5	0.97	0.05	74,86,87,89	0
2	PO4	A	301	5/5	0.97	0.05	65,67,76,79	0
2	PO4	F	301	5/5	0.98	0.04	47,57,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.