Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:43 PM GMT

PDB ID	:	1GXB
Title	:	ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE IN COMPLEX WITH PYROPHOSPHATE AND MAGNESIUM
Authors	:	Mayans, O.; Ivens, A.; Nissen, L.J.; Kirschner, K.; Wilmanns, M.
Deposited on	:	2002-04-02
Resolution	:	2.65 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

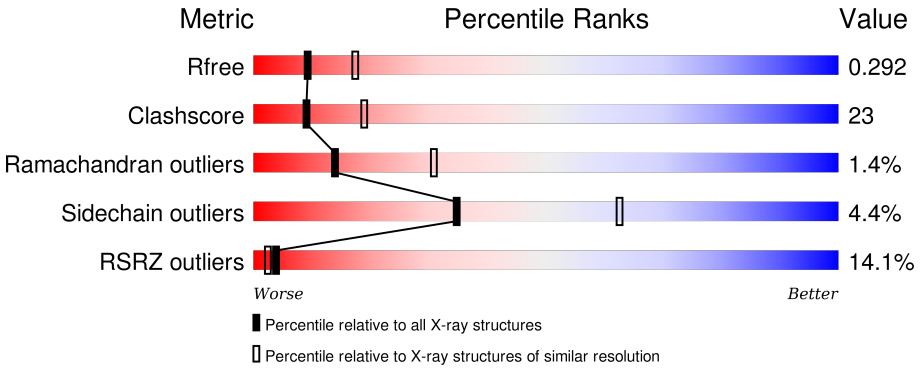
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3152 (2.70-2.62)
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)
RSRZ outliers	91569	3161 (2.70-2.62)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	345	15% 63% 32%
1	B	345	12% 58% 35% 5%
1	C	345	10% 55% 39%
1	D	345	19% 60% 37%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	POP	D	1349	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10753 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.

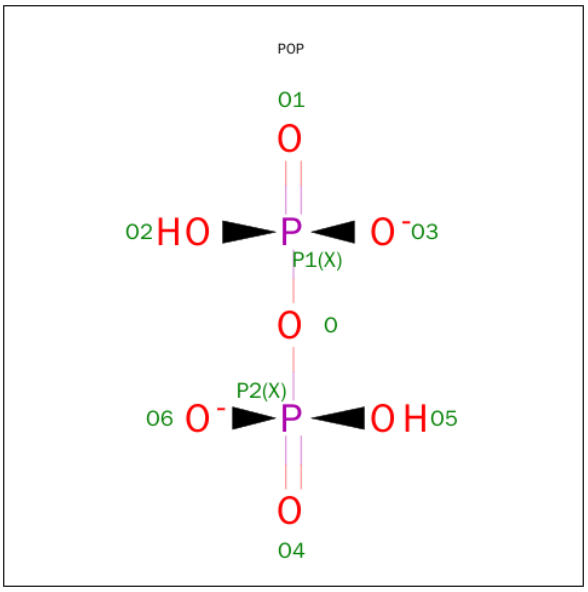
- Molecule 1 is a protein called ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	1
			2609	1674	440	488	7			
1	B	340	Total	C	N	O	S	0	0	1
			2611	1675	440	489	7			
1	C	341	Total	C	N	O	S	0	0	0
			2619	1679	441	492	7			
1	D	345	Total	C	N	O	S	0	0	0
			2642	1691	445	499	7			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: $H_2O_7P_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 9 7 2	0	0
3	B	1	Total O P 9 7 2	0	0
3	C	1	Total O P 9 7 2	0	0
3	D	1	Total O P 9 7 2	0	0

- Molecule 4 is water.

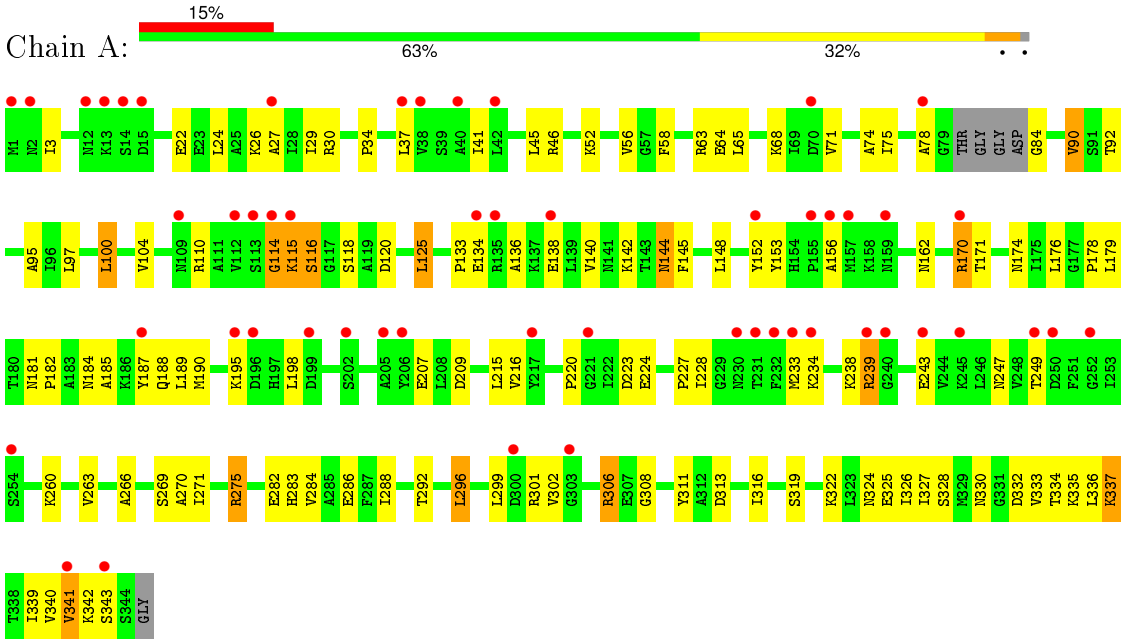
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	61	Total O 61 61	0	0
4	B	56	Total O 56 56	0	0
4	C	66	Total O 66 66	0	0
4	D	50	Total O 50 50	0	0

3 Residue-property plots (i)

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

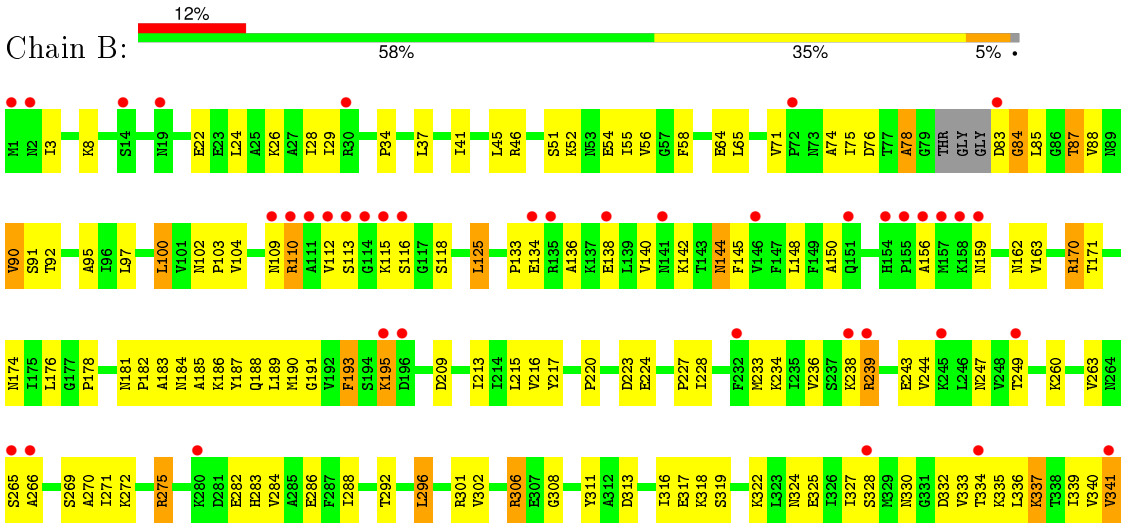
- Molecule 1: ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE

Chain A:



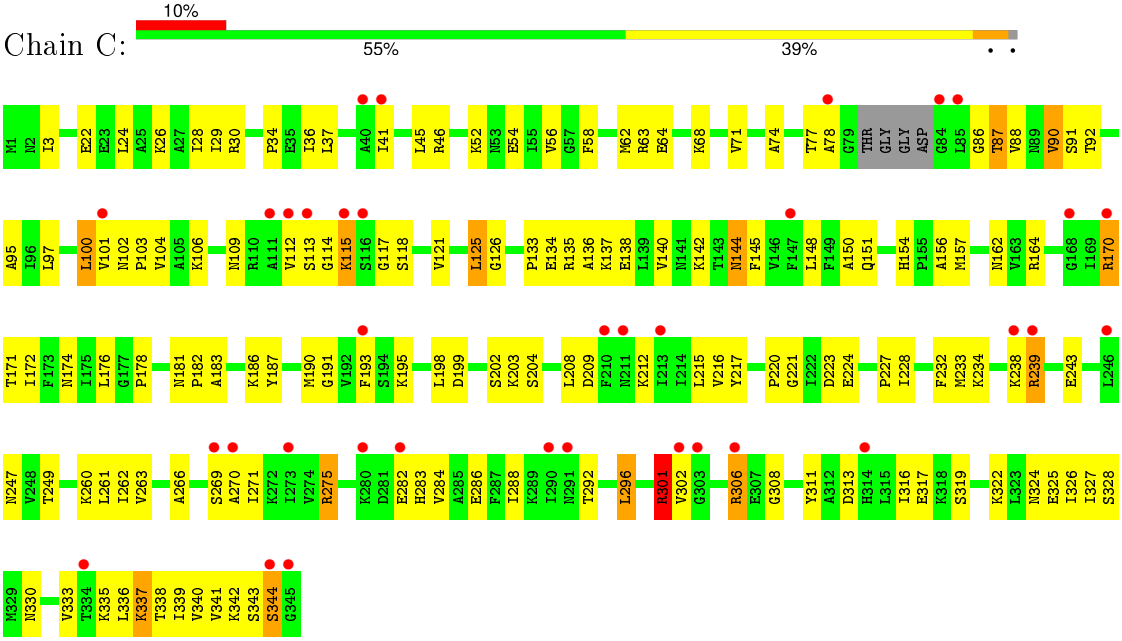
- Molecule 1: ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE

Chain B:



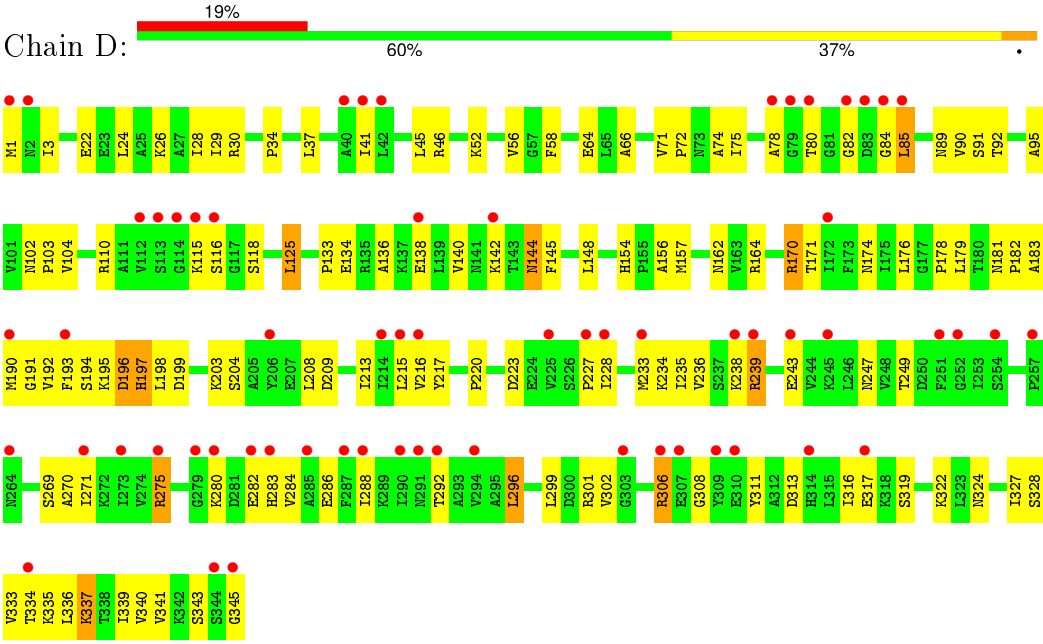
• Molecule 1: ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE

Chain C:



• Molecule 1: ANTHRANILATE PHOSPHORIBOSYLTRANSFERASE

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	90.84Å 65.47Å 115.37Å 90.00° 107.25° 90.00°	Depositor
Resolution (Å)	20.00 – 2.65 19.69 – 2.63	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.65) 99.0 (19.69-2.63)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.88 (at 2.63Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.223 , 0.297 0.282 , 0.292	Depositor DCC
R_{free} test set	1141 reflections (3.02%)	DCC
Wilson B-factor (Å^2)	62.7	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 64.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 38448 reflections	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	10753	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 20.66 % 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 8.2807e-03.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, POP

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	0/2645	0.97	11/3573 (0.3%)
1	B	0.38	0/2647	0.83	10/3576 (0.3%)
1	C	0.40	0/2655	0.99	11/3584 (0.3%)
1	D	0.36	0/2679	0.82	10/3618 (0.3%)
All	All	0.39	0/10626	0.91	42/14351 (0.3%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	301	ARG	NE-CZ-NH1	-23.00	108.80	120.30
1	C	301	ARG	NE-CZ-NH2	20.39	130.49	120.30
1	A	306	ARG	NE-CZ-NH2	19.56	130.08	120.30
1	A	306	ARG	NE-CZ-NH1	-19.42	110.59	120.30
1	A	275	ARG	NE-CZ-NH1	-15.11	112.75	120.30
1	D	275	ARG	NE-CZ-NH2	-15.10	112.75	120.30
1	C	275	ARG	NE-CZ-NH1	-14.95	112.83	120.30
1	B	275	ARG	NE-CZ-NH2	-14.85	112.88	120.30
1	A	275	ARG	NE-CZ-NH2	14.35	127.47	120.30
1	C	275	ARG	NE-CZ-NH2	14.23	127.42	120.30
1	D	275	ARG	NE-CZ-NH1	14.14	127.37	120.30
1	B	275	ARG	NE-CZ-NH1	13.98	127.29	120.30
1	B	170	ARG	NE-CZ-NH2	-13.79	113.41	120.30
1	C	170	ARG	NE-CZ-NH1	-13.61	113.50	120.30
1	A	170	ARG	NE-CZ-NH2	-13.60	113.50	120.30
1	C	170	ARG	NE-CZ-NH2	13.34	126.97	120.30
1	D	170	ARG	NE-CZ-NH1	-13.01	113.80	120.30
1	B	170	ARG	NE-CZ-NH1	12.93	126.77	120.30
1	A	170	ARG	NE-CZ-NH1	12.72	126.66	120.30
1	D	170	ARG	NE-CZ-NH2	12.38	126.49	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	301	ARG	CD-NE-CZ	10.54	138.36	123.60
1	A	306	ARG	CD-NE-CZ	9.51	136.92	123.60
1	D	301	ARG	NE-CZ-NH2	-7.89	116.35	120.30
1	B	275	ARG	CD-NE-CZ	7.63	134.28	123.60
1	D	275	ARG	CD-NE-CZ	7.61	134.26	123.60
1	C	275	ARG	CD-NE-CZ	7.59	134.22	123.60
1	A	301	ARG	NE-CZ-NH2	-7.53	116.54	120.30
1	A	275	ARG	CD-NE-CZ	7.51	134.11	123.60
1	A	301	ARG	NE-CZ-NH1	7.51	124.05	120.30
1	D	301	ARG	NE-CZ-NH1	6.85	123.73	120.30
1	C	170	ARG	CD-NE-CZ	6.76	133.06	123.60
1	C	306	ARG	NE-CZ-NH2	-6.65	116.97	120.30
1	D	306	ARG	NE-CZ-NH2	-6.65	116.98	120.30
1	B	306	ARG	NE-CZ-NH2	-6.63	116.98	120.30
1	D	170	ARG	CD-NE-CZ	6.54	132.76	123.60
1	B	306	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	B	301	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	170	ARG	CD-NE-CZ	6.38	132.54	123.60
1	B	301	ARG	NE-CZ-NH1	6.38	123.49	120.30
1	B	170	ARG	CD-NE-CZ	6.37	132.51	123.60
1	D	306	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	C	306	ARG	NE-CZ-NH1	6.24	123.42	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2609	0	2725	109	0
1	B	2611	0	2725	128	0
1	C	2619	0	2734	127	0
1	D	2642	0	2752	130	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	9	0	0	0	0
3	B	9	0	0	2	0
3	C	9	0	0	1	0
3	D	9	0	0	4	0
4	A	61	0	0	11	0
4	B	56	0	0	14	0
4	C	66	0	0	19	0
4	D	50	0	0	23	0
All	All	10753	0	10936	496	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 23.

All (496) 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:C:109:ASN:HD22	1:C:150:ALA:HB3	1.10	1.12
1:A:78:ALA:HB2	1:A:188:GLN:OE1	1.62	0.98
1:D:280:LYS:HE3	1:D:345:GLY:H	1.32	0.93
1:A:178:PRO:HB2	1:A:190:MET:HE2	1.48	0.93
1:C:171:THR:H	1:C:174:ASN:ND2	1.67	0.92
1:D:171:THR:H	1:D:174:ASN:ND2	1.66	0.92
1:A:171:THR:H	1:A:174:ASN:ND2	1.70	0.90
1:B:171:THR:H	1:B:174:ASN:ND2	1.70	0.89
1:A:84:GLY:N	4:A:2016:HOH:O	2.06	0.89
1:B:83:ASP:HB3	4:B:2014:HOH:O	1.73	0.88
1:B:78:ALA:HB3	1:B:190:MET:HA	1.53	0.88
1:B:78:ALA:HB2	1:B:188:GLN:NE2	1.88	0.88
1:D:82:GLY:CA	1:D:170:ARG:HH22	1.87	0.86
1:D:235:ILE:HD11	4:D:2036:HOH:O	1.76	0.85
1:C:342:LYS:HD2	4:C:2055:HOH:O	1.75	0.85
1:A:29:ILE:CG2	1:A:65:LEU:HD12	2.08	0.83
1:A:78:ALA:HA	4:A:2018:HOH:O	1.78	0.82
1:D:171:THR:H	1:D:174:ASN:HD22	1.25	0.82
1:C:109:ASN:ND2	1:C:150:ALA:HB3	1.94	0.81
1:C:62:MET:HE1	4:C:2017:HOH:O	1.80	0.81
1:C:109:ASN:HD21	1:C:151:GLN:HG3	1.44	0.81
1:C:88:VAL:HG13	1:C:261:LEU:O	1.80	0.81
1:A:116:SER:HA	4:A:2044:HOH:O	1.81	0.80
1:D:208:LEU:HG	4:D:2013:HOH:O	1.81	0.80
1:B:171:THR:H	1:B:174:ASN:HD22	1.28	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:ALA:CB	1:D:178:PRO:HB3	2.11	0.80
1:C:171:THR:H	1:C:174:ASN:HD22	1.28	0.79
1:D:181:ASN:HD21	1:D:183:ALA:HB3	1.47	0.79
1:A:178:PRO:HB2	1:A:190:MET:CE	2.13	0.78
1:A:171:THR:H	1:A:174:ASN:HD22	1.29	0.78
1:A:266:ALA:HA	4:A:2044:HOH:O	1.84	0.78
1:D:170:ARG:HB3	4:D:2027:HOH:O	1.82	0.77
1:B:188:GLN:NE2	1:B:190:MET:HB2	1.99	0.77
1:B:181:ASN:ND2	1:B:185:ALA:H	1.82	0.77
1:C:97:LEU:HD22	1:C:319:SER:OG	1.85	0.77
1:B:97:LEU:HD22	1:B:319:SER:OG	1.85	0.76
1:D:82:GLY:HA2	1:D:170:ARG:HH22	1.47	0.76
1:D:190:MET:O	1:D:215:LEU:HA	1.85	0.76
1:A:97:LEU:HD22	1:A:319:SER:OG	1.86	0.75
1:D:97:LEU:HD22	1:D:319:SER:OG	1.85	0.75
1:B:78:ALA:CB	1:B:188:GLN:NE2	2.50	0.74
1:D:164:ARG:NH1	4:D:2025:HOH:O	2.19	0.74
1:B:87:THR:HG23	1:B:88:VAL:O	1.86	0.74
1:C:78:ALA:CB	1:C:178:PRO:HB3	2.17	0.74
1:D:216:VAL:HG12	1:D:233:MET:HB3	1.69	0.74
1:B:234:LYS:HE2	1:B:243:GLU:HG3	1.70	0.74
1:B:327:ILE:HD12	1:B:336:LEU:HD22	1.69	0.74
1:C:234:LYS:HE2	1:C:243:GLU:HG3	1.70	0.73
1:D:235:ILE:CG1	4:D:2036:HOH:O	2.36	0.73
1:D:234:LYS:HE2	1:D:243:GLU:HG3	1.71	0.73
1:D:82:GLY:HA2	1:D:170:ARG:NH2	2.04	0.72
1:C:216:VAL:HG12	1:C:233:MET:HB3	1.69	0.72
1:A:29:ILE:HG21	1:A:65:LEU:HD12	1.71	0.72
1:C:88:VAL:HG12	1:C:263:VAL:HG13	1.70	0.72
1:D:78:ALA:HB1	1:D:178:PRO:HB3	1.72	0.72
1:D:328:SER:HA	1:D:333:VAL:HG22	1.73	0.71
1:A:234:LYS:HE2	1:A:243:GLU:HG3	1.73	0.71
1:A:216:VAL:HG12	1:A:233:MET:HB3	1.70	0.71
1:B:148:LEU:HD13	1:B:182:PRO:HB2	1.71	0.71
1:A:336:LEU:O	1:A:340:VAL:HG23	1.90	0.71
1:D:327:ILE:HD12	1:D:336:LEU:HD22	1.71	0.70
1:C:327:ILE:HD12	1:C:336:LEU:HD22	1.72	0.70
1:A:327:ILE:HD12	1:A:336:LEU:HD22	1.74	0.70
1:A:75:ILE:HA	1:A:187:TYR:O	1.91	0.70
1:C:328:SER:HA	1:C:333:VAL:HG22	1.74	0.69
1:B:193:PHE:HD1	1:B:193:PHE:H	1.40	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:328:SER:HA	1:B:333:VAL:HG22	1.73	0.69
1:D:239:ARG:HD3	1:D:239:ARG:H	1.59	0.68
1:D:171:THR:N	1:D:174:ASN:HD22	1.91	0.68
1:A:328:SER:HA	1:A:333:VAL:HG22	1.74	0.68
1:C:28:ILE:HA	4:C:2008:HOH:O	1.93	0.68
1:B:324:ASN:ND2	4:B:2052:HOH:O	2.26	0.68
1:A:239:ARG:H	1:A:239:ARG:HD3	1.57	0.67
1:D:170:ARG:HG2	4:D:2010:HOH:O	1.95	0.67
1:A:29:ILE:HG22	1:A:65:LEU:HD12	1.76	0.67
1:C:336:LEU:O	1:C:340:VAL:HG23	1.94	0.67
1:B:216:VAL:HG12	1:B:233:MET:HB3	1.75	0.67
1:D:336:LEU:O	1:D:340:VAL:HG23	1.93	0.67
1:B:116:SER:OG	4:B:2021:HOH:O	2.13	0.67
1:C:137:LYS:HB2	4:C:2027:HOH:O	1.95	0.66
1:D:275:ARG:HG2	1:D:343:SER:HB2	1.77	0.66
1:C:171:THR:N	1:C:174:ASN:HD22	1.93	0.66
1:C:239:ARG:HD3	1:C:239:ARG:H	1.59	0.66
1:D:80:THR:HG21	1:D:193:PHE:H	1.61	0.66
1:C:172:ILE:HA	4:C:2041:HOH:O	1.95	0.66
1:B:336:LEU:O	1:B:340:VAL:HG23	1.96	0.65
1:A:156:ALA:HB2	4:A:2013:HOH:O	1.96	0.65
1:B:171:THR:N	1:B:174:ASN:HD22	1.95	0.65
1:D:116:SER:O	4:D:2019:HOH:O	2.15	0.65
1:D:233:MET:SD	4:D:2036:HOH:O	2.54	0.65
1:B:181:ASN:HD22	1:B:185:ALA:H	1.43	0.64
1:A:125:LEU:HB3	1:A:327:ILE:HD11	1.79	0.64
1:B:239:ARG:HD3	1:B:239:ARG:H	1.61	0.64
3:D:1349:POP:P1	4:D:2048:HOH:O	2.54	0.64
1:B:324:ASN:HD21	1:B:337:LYS:HE3	1.63	0.64
1:C:125:LEU:HB3	1:C:327:ILE:HD11	1.78	0.64
1:B:125:LEU:HB3	1:B:327:ILE:HD11	1.80	0.64
1:D:333:VAL:O	1:D:336:LEU:HB3	1.97	0.64
1:D:125:LEU:HB3	1:D:327:ILE:HD11	1.79	0.64
1:C:125:LEU:O	1:C:327:ILE:HD13	1.98	0.64
1:B:333:VAL:O	1:B:336:LEU:HB3	1.98	0.63
1:B:85:LEU:HD23	1:B:85:LEU:O	1.97	0.63
1:A:134:GLU:HG3	1:D:328:SER:O	1.98	0.63
1:A:171:THR:N	1:A:174:ASN:HD22	1.95	0.63
1:C:324:ASN:HD21	1:C:337:LYS:HE3	1.63	0.63
1:A:333:VAL:O	1:A:336:LEU:HB3	1.99	0.63
1:D:78:ALA:HB3	1:D:178:PRO:HB3	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:LYS:HG2	1:C:266:ALA:HB2	1.80	0.63
1:A:223:ASP:OD1	4:A:2034:HOH:O	2.15	0.63
1:B:340:VAL:C	1:B:342:LYS:H	2.02	0.62
1:D:125:LEU:O	1:D:327:ILE:HD13	2.00	0.62
1:A:148:LEU:HD13	1:A:182:PRO:HB2	1.81	0.62
1:B:271:ILE:O	1:B:275:ARG:HG3	1.99	0.62
1:D:46:ARG:HD2	1:D:170:ARG:O	2.00	0.61
1:B:76:ASP:HB3	1:B:188:GLN:HG3	1.81	0.61
1:C:46:ARG:HD2	1:C:170:ARG:O	2.01	0.61
1:C:333:VAL:O	1:C:336:LEU:HB3	2.01	0.61
3:D:1349:POP:O1	4:D:2048:HOH:O	2.15	0.61
1:A:125:LEU:O	1:A:327:ILE:HD13	2.01	0.61
1:A:333:VAL:HG23	4:A:2058:HOH:O	2.01	0.61
1:A:335:LYS:HG3	4:A:2059:HOH:O	2.00	0.61
1:D:324:ASN:HD21	1:D:337:LYS:HE3	1.63	0.60
1:D:24:LEU:HD11	1:D:41:ILE:HD13	1.82	0.60
1:D:181:ASN:ND2	1:D:183:ALA:H	2.00	0.60
1:A:271:ILE:O	1:A:275:ARG:HG3	2.02	0.60
1:D:82:GLY:C	1:D:170:ARG:HH22	2.04	0.60
1:A:34:PRO:HD2	1:A:37:LEU:HD12	1.84	0.60
1:B:78:ALA:HB1	1:B:178:PRO:HB3	1.84	0.60
1:D:271:ILE:O	1:D:275:ARG:HG3	2.01	0.59
1:B:34:PRO:HD2	1:B:37:LEU:HD12	1.85	0.59
1:A:324:ASN:HD21	1:A:337:LYS:HE3	1.67	0.59
1:C:24:LEU:HD11	1:C:41:ILE:HD13	1.84	0.59
1:B:188:GLN:HE22	1:B:190:MET:HB2	1.63	0.59
1:B:193:PHE:CD1	1:B:193:PHE:N	2.69	0.59
1:B:46:ARG:HD2	1:B:170:ARG:O	2.03	0.59
1:C:88:VAL:HG12	1:C:263:VAL:CG1	2.33	0.58
1:C:109:ASN:ND2	1:C:151:GLN:HG3	2.16	0.58
1:C:68:LYS:HE3	4:C:2016:HOH:O	2.03	0.58
1:C:199:ASP:OD2	1:C:203:LYS:HE3	2.04	0.58
1:B:29:ILE:CG2	1:B:65:LEU:HD12	2.33	0.58
1:A:46:ARG:HD2	1:A:170:ARG:O	2.03	0.58
1:A:195:LYS:O	1:A:198:LEU:HB2	2.04	0.58
1:D:34:PRO:HD2	1:D:37:LEU:HD12	1.85	0.58
1:A:340:VAL:O	1:A:342:LYS:N	2.37	0.58
1:B:125:LEU:O	1:B:327:ILE:HD13	2.04	0.57
1:B:75:ILE:HA	1:B:187:TYR:O	2.05	0.57
1:A:78:ALA:HB3	1:A:190:MET:HA	1.86	0.57
1:B:24:LEU:HD11	1:B:41:ILE:HD13	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:VAL:CG2	4:C:2057:HOH:O	2.51	0.57
1:A:27:ALA:HB1	4:A:2005:HOH:O	2.05	0.57
1:C:34:PRO:HD2	1:C:37:LEU:HD12	1.87	0.57
1:B:178:PRO:HB2	1:B:190:MET:SD	2.45	0.56
1:B:184:ASN:ND2	1:B:186:LYS:HE2	2.20	0.56
1:B:341:VAL:HG12	1:B:341:VAL:O	2.04	0.56
1:D:235:ILE:CD1	4:D:2036:HOH:O	2.40	0.56
1:D:28:ILE:HA	4:D:2006:HOH:O	2.05	0.56
1:C:216:VAL:HG12	1:C:233:MET:CB	2.36	0.56
1:A:216:VAL:HG12	1:A:233:MET:CB	2.36	0.56
1:C:26:LYS:HE2	1:C:64:GLU:OE1	2.06	0.56
1:D:216:VAL:HG12	1:D:233:MET:CB	2.34	0.56
1:B:41:ILE:O	1:B:45:LEU:HB2	2.06	0.56
1:A:190:MET:O	1:A:215:LEU:HA	2.06	0.55
1:B:195:LYS:HG2	1:B:217:TYR:CD2	2.42	0.55
1:C:186:LYS:HD2	4:C:2046:HOH:O	2.06	0.55
1:D:26:LYS:HE2	1:D:64:GLU:OE1	2.05	0.55
1:C:109:ASN:HD21	1:C:151:GLN:CG	2.18	0.55
1:C:36:ILE:HG13	4:C:2010:HOH:O	2.06	0.55
1:A:24:LEU:HD11	1:A:41:ILE:HD13	1.88	0.55
1:C:121:VAL:HG21	4:C:2057:HOH:O	2.06	0.55
1:C:78:ALA:HB1	1:C:178:PRO:HB3	1.86	0.54
1:B:125:LEU:CD1	1:B:270:ALA:HB1	2.38	0.54
1:C:126:GLY:HA2	4:C:2024:HOH:O	2.07	0.54
1:B:26:LYS:HE2	1:B:64:GLU:OE1	2.07	0.54
1:A:92:THR:OG1	1:A:118:SER:HB2	2.08	0.54
1:A:78:ALA:CB	1:A:188:GLN:OE1	2.48	0.54
1:B:22:GLU:HG2	1:B:26:LYS:HE3	1.90	0.54
1:A:134:GLU:H	1:A:134:GLU:CD	2.12	0.53
1:A:26:LYS:HE2	1:A:64:GLU:OE1	2.07	0.53
1:D:220:PRO:HD3	1:D:228:ILE:HD13	1.91	0.53
1:C:271:ILE:O	1:C:275:ARG:HG3	2.07	0.53
1:D:92:THR:CB	4:D:2016:HOH:O	2.55	0.53
1:D:170:ARG:HA	1:D:174:ASN:ND2	2.23	0.53
1:D:134:GLU:CD	1:D:134:GLU:H	2.12	0.53
1:D:148:LEU:HD13	1:D:182:PRO:HB2	1.90	0.53
1:B:178:PRO:HB2	1:B:190:MET:HG3	1.91	0.53
1:B:216:VAL:HG12	1:B:233:MET:CB	2.39	0.53
1:B:92:THR:OG1	1:B:118:SER:HB2	2.09	0.53
1:D:191:GLY:O	1:D:223:ASP:HB2	2.09	0.53
1:C:125:LEU:CD1	1:C:270:ALA:HB1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:22:GLU:HG2	1:C:26:LYS:HE3	1.92	0.52
1:A:292:THR:HG22	1:A:296:LEU:HD22	1.91	0.52
1:C:171:THR:N	1:C:174:ASN:ND2	2.47	0.52
1:D:118:SER:HB3	4:D:2020:HOH:O	2.08	0.52
1:D:41:ILE:O	1:D:45:LEU:HB2	2.09	0.52
1:A:95:ALA:HB1	1:A:104:VAL:HG11	1.92	0.52
1:D:194:SER:HA	1:D:217:TYR:CE1	2.45	0.52
1:B:340:VAL:O	1:B:342:LYS:N	2.43	0.52
1:D:85:LEU:N	1:D:85:LEU:HD22	2.24	0.52
1:C:170:ARG:HA	1:C:174:ASN:ND2	2.23	0.52
1:C:343:SER:O	1:C:344:SER:CB	2.57	0.52
1:D:3:ILE:HG23	1:D:41:ILE:HD11	1.92	0.52
1:A:22:GLU:HG2	1:A:26:LYS:HE3	1.91	0.52
1:D:92:THR:OG1	1:D:118:SER:HB2	2.10	0.52
1:D:136:ALA:O	1:D:140:VAL:HG23	2.10	0.52
1:B:292:THR:HG22	1:B:296:LEU:HD22	1.91	0.52
1:B:134:GLU:CD	1:B:134:GLU:H	2.13	0.52
1:C:150:ALA:O	1:C:154:HIS:HB2	2.10	0.52
1:D:181:ASN:ND2	1:D:183:ALA:HB3	2.21	0.52
1:C:193:PHE:O	1:C:217:TYR:HE1	1.93	0.52
1:A:125:LEU:CD1	1:A:270:ALA:HB1	2.40	0.52
1:A:341:VAL:C	1:A:343:SER:H	2.12	0.52
1:B:75:ILE:HG13	1:B:187:TYR:O	2.10	0.52
1:B:318:LYS:NZ	4:B:2049:HOH:O	2.43	0.52
1:C:342:LYS:C	1:C:344:SER:H	2.13	0.51
1:D:22:GLU:HG2	1:D:26:LYS:HE3	1.91	0.51
1:B:220:PRO:HD3	1:B:228:ILE:HD13	1.91	0.51
1:A:170:ARG:HA	1:A:174:ASN:ND2	2.25	0.51
1:C:292:THR:HG22	1:C:296:LEU:HD22	1.91	0.51
1:C:3:ILE:HG23	1:C:41:ILE:HD11	1.92	0.51
1:C:187:TYR:N	1:C:187:TYR:CD1	2.78	0.51
1:C:134:GLU:CD	1:C:134:GLU:H	2.12	0.51
1:D:233:MET:HG3	4:D:2036:HOH:O	2.09	0.51
1:A:41:ILE:O	1:A:45:LEU:HB2	2.10	0.51
1:A:136:ALA:O	1:A:140:VAL:HG23	2.11	0.51
1:C:91:SER:HB2	3:C:1348:POP:P2	2.51	0.51
1:B:115:LYS:HE3	1:B:266:ALA:HB2	1.93	0.51
1:D:91:SER:HB2	3:D:1349:POP:O	2.11	0.50
1:A:114:GLY:O	1:A:115:LYS:C	2.50	0.50
1:D:80:THR:HG21	1:D:192:VAL:HA	1.93	0.50
1:D:292:THR:HG22	1:D:296:LEU:HD22	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:ARG:HA	1:B:174:ASN:ND2	2.26	0.50
1:D:140:VAL:O	1:D:144:ASN:HA	2.12	0.50
1:A:341:VAL:HG12	1:A:341:VAL:O	2.11	0.50
1:C:133:PRO:HG2	1:C:134:GLU:OE2	2.11	0.50
1:C:136:ALA:O	1:C:140:VAL:HG23	2.11	0.50
1:D:80:THR:CG2	1:D:192:VAL:HA	2.41	0.50
1:D:110:ARG:HG2	1:D:110:ARG:HH11	1.75	0.50
1:B:78:ALA:HB3	1:B:190:MET:CA	2.35	0.50
1:D:125:LEU:CD1	1:D:270:ALA:HB1	2.42	0.50
1:C:335:LYS:O	1:C:339:ILE:HG12	2.12	0.50
1:D:263:VAL:HG11	1:D:269:SER:HA	1.94	0.50
1:A:152:TYR:HD2	1:A:153:TYR:CE1	2.30	0.50
1:D:239:ARG:HG3	1:D:239:ARG:HH11	1.77	0.49
1:D:80:THR:HG21	1:D:193:PHE:N	2.27	0.49
1:B:159:ASN:HA	4:B:2026:HOH:O	2.10	0.49
1:C:190:MET:O	1:C:215:LEU:HA	2.11	0.49
1:C:112:VAL:O	1:C:114:GLY:N	2.45	0.49
1:B:136:ALA:O	1:B:140:VAL:HG23	2.12	0.49
1:B:140:VAL:O	1:B:144:ASN:HA	2.11	0.49
1:B:112:VAL:HG12	1:B:112:VAL:O	2.12	0.49
1:D:260:LYS:HG2	1:D:283:HIS:CD2	2.47	0.49
1:A:335:LYS:O	1:A:339:ILE:HG12	2.12	0.49
1:A:260:LYS:HG2	1:A:283:HIS:CD2	2.47	0.49
1:A:247:ASN:ND2	1:A:249:THR:HG23	2.27	0.49
1:C:140:VAL:O	1:C:144:ASN:HA	2.12	0.49
1:B:335:LYS:O	1:B:339:ILE:HG12	2.12	0.49
1:B:247:ASN:ND2	1:B:249:THR:HG23	2.27	0.49
1:B:263:VAL:HG11	1:B:269:SER:HA	1.93	0.49
1:B:95:ALA:HB1	1:B:104:VAL:HG11	1.94	0.49
1:B:3:ILE:HG23	1:B:41:ILE:HD11	1.94	0.49
1:C:112:VAL:HG11	4:C:2015:HOH:O	2.11	0.49
1:D:247:ASN:ND2	1:D:249:THR:HG23	2.28	0.49
1:C:220:PRO:HD3	1:C:228:ILE:HD13	1.94	0.49
1:B:272:LYS:NZ	4:B:2040:HOH:O	2.45	0.49
1:B:125:LEU:HD13	1:B:270:ALA:HB1	1.95	0.49
1:D:195:LYS:O	1:D:198:LEU:HB3	2.13	0.49
1:D:302:VAL:HG11	1:D:308:GLY:HA2	1.94	0.49
1:D:335:LYS:O	1:D:339:ILE:HG12	2.12	0.49
1:B:260:LYS:HG2	1:B:283:HIS:CD2	2.48	0.49
1:C:313:ASP:O	1:C:316:ILE:HG22	2.12	0.49
1:C:125:LEU:HD13	1:C:270:ALA:HB1	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:260:LYS:HG2	1:C:283:HIS:CD2	2.47	0.48
1:A:239:ARG:HH11	1:A:239:ARG:HG3	1.78	0.48
1:C:41:ILE:O	1:C:45:LEU:HB2	2.14	0.48
1:C:157:MET:HE1	4:C:2043:HOH:O	2.13	0.48
1:A:63:ARG:O	1:A:68:LYS:HE3	2.13	0.48
1:C:148:LEU:HD13	1:C:182:PRO:HB2	1.95	0.48
1:C:239:ARG:HG3	1:C:239:ARG:HH11	1.77	0.48
1:B:156:ALA:HB3	4:B:2025:HOH:O	2.12	0.48
1:B:29:ILE:HG21	1:B:65:LEU:HD12	1.95	0.48
1:A:302:VAL:HG11	1:A:308:GLY:HA2	1.95	0.48
1:D:271:ILE:HD12	4:D:2041:HOH:O	2.12	0.48
1:D:195:LYS:HG3	1:D:217:TYR:CE2	2.48	0.48
1:D:66:ALA:HB2	1:D:154:HIS:CE1	2.48	0.48
1:C:263:VAL:HG11	1:C:269:SER:HA	1.95	0.48
1:A:3:ILE:HG23	1:A:41:ILE:HD11	1.95	0.48
1:A:220:PRO:HD3	1:A:228:ILE:HD13	1.95	0.48
1:A:138:GLU:OE2	1:A:142:LYS:HE2	2.14	0.48
1:C:302:VAL:HG11	1:C:308:GLY:HA2	1.95	0.48
1:A:133:PRO:HG2	1:A:134:GLU:OE2	2.14	0.48
1:C:247:ASN:ND2	1:C:249:THR:HG23	2.29	0.48
1:C:101:VAL:O	1:C:301:ARG:NH1	2.47	0.48
1:C:342:LYS:C	1:C:344:SER:N	2.66	0.48
1:D:235:ILE:HG13	4:D:2036:HOH:O	2.09	0.48
1:D:100:LEU:HD13	1:D:322:LYS:HG2	1.97	0.47
1:B:138:GLU:OE2	1:B:142:LYS:HE2	2.14	0.47
1:D:78:ALA:O	1:D:190:MET:HA	2.14	0.47
1:C:227:PRO:HG2	1:C:228:ILE:HG23	1.97	0.47
1:C:191:GLY:HA3	1:C:223:ASP:O	2.14	0.47
1:C:164:ARG:HD2	4:C:2040:HOH:O	2.14	0.47
1:B:109:ASN:ND2	1:B:150:ALA:HB3	2.29	0.47
1:B:8:LYS:NZ	4:B:2002:HOH:O	2.47	0.47
1:C:181:ASN:HD21	1:C:183:ALA:HB3	1.79	0.47
1:A:140:VAL:O	1:A:144:ASN:HA	2.14	0.47
1:C:284:VAL:O	1:C:288:ILE:HG13	2.15	0.47
1:C:170:ARG:HG2	4:C:2011:HOH:O	2.14	0.47
1:C:78:ALA:HB3	1:C:178:PRO:HB3	1.97	0.47
1:B:302:VAL:HG11	1:B:308:GLY:HA2	1.97	0.47
1:A:263:VAL:HG11	1:A:269:SER:HA	1.96	0.47
1:C:92:THR:OG1	1:C:118:SER:HB2	2.14	0.46
1:C:204:SER:O	1:C:208:LEU:HD23	2.14	0.46
1:C:209:ASP:OD1	1:C:238:LYS:NZ	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:199:ASP:OD1	1:D:203:LYS:HE3	2.16	0.46
1:B:328:SER:HA	1:B:333:VAL:CG2	2.43	0.46
1:C:198:LEU:HD11	1:C:232:PHE:O	2.15	0.46
1:B:239:ARG:HH11	1:B:239:ARG:HG3	1.80	0.46
1:B:341:VAL:C	1:B:342:LYS:HE3	2.36	0.46
1:D:24:LEU:HD11	1:D:41:ILE:CD1	2.45	0.46
1:D:284:VAL:O	1:D:288:ILE:HG13	2.15	0.46
3:B:1347:POP:O	4:B:2056:HOH:O	2.21	0.46
1:C:138:GLU:OE2	1:C:142:LYS:HE2	2.15	0.46
1:A:328:SER:HA	1:A:333:VAL:CG2	2.44	0.46
1:C:328:SER:HA	1:C:333:VAL:CG2	2.44	0.46
1:B:133:PRO:HG2	1:B:134:GLU:OE2	2.15	0.46
1:A:52:LYS:O	1:A:56:VAL:HG23	2.15	0.46
1:B:195:LYS:HG2	1:B:217:TYR:CE2	2.51	0.46
1:C:95:ALA:HB1	1:C:104:VAL:HG11	1.98	0.46
1:D:179:LEU:HD22	4:D:2013:HOH:O	2.14	0.46
1:C:198:LEU:HD22	1:C:234:LYS:HG2	1.98	0.46
1:B:184:ASN:HD21	1:B:186:LYS:HE2	1.80	0.46
1:A:22:GLU:O	1:A:26:LYS:HG3	2.16	0.46
1:D:71:VAL:HG12	1:D:74:ALA:HB2	1.97	0.46
1:B:342:LYS:HE3	1:B:342:LYS:N	2.31	0.46
1:A:110:ARG:NH1	1:A:120:ASP:OD1	2.46	0.46
1:C:90:VAL:HG21	1:C:224:GLU:HG2	1.98	0.46
1:D:171:THR:N	1:D:174:ASN:ND2	2.47	0.45
1:A:187:TYR:CD2	1:A:299:LEU:HD23	2.51	0.45
1:C:30:ARG:NH2	4:C:2007:HOH:O	2.25	0.45
1:A:313:ASP:O	1:A:316:ILE:HG22	2.15	0.45
1:A:115:LYS:HD3	1:A:266:ALA:CB	2.46	0.45
1:D:186:LYS:HD3	1:D:186:LYS:HA	1.69	0.45
1:B:340:VAL:C	1:B:342:LYS:N	2.69	0.45
1:A:125:LEU:HD13	1:A:270:ALA:HB1	1.97	0.45
1:D:220:PRO:HD3	1:D:228:ILE:CD1	2.46	0.45
1:B:91:SER:HB2	3:B:1347:POP:O2	2.17	0.45
1:A:138:GLU:O	1:A:142:LYS:HG2	2.16	0.45
1:D:145:PHE:C	1:D:145:PHE:CD2	2.89	0.45
1:C:86:GLY:O	1:C:262:ILE:HG23	2.17	0.45
1:A:171:THR:N	1:A:174:ASN:ND2	2.50	0.45
1:D:328:SER:HA	1:D:333:VAL:CG2	2.43	0.45
1:D:91:SER:HB2	3:D:1349:POP:P2	2.56	0.45
1:D:92:THR:HB	4:D:2016:HOH:O	2.16	0.45
1:D:196:ASP:O	1:D:198:LEU:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:LYS:NZ	1:A:207:GLU:OE2	2.45	0.45
1:D:138:GLU:OE2	1:D:142:LYS:HE2	2.17	0.45
1:B:113:SER:HA	4:B:2020:HOH:O	2.16	0.45
1:B:209:ASP:OD1	1:B:238:LYS:NZ	2.50	0.45
1:A:341:VAL:C	4:A:2060:HOH:O	2.54	0.45
1:D:227:PRO:HG2	1:D:228:ILE:HG23	1.98	0.45
1:A:144:ASN:HD22	1:A:144:ASN:HA	1.61	0.45
1:A:209:ASP:OD1	1:A:238:LYS:NZ	2.50	0.45
1:A:90:VAL:HG21	1:A:224:GLU:HG2	1.99	0.45
1:A:29:ILE:HG22	1:A:156:ALA:HB1	1.99	0.45
1:B:227:PRO:HG2	1:B:228:ILE:HG23	1.99	0.45
1:C:77:THR:O	1:C:106:LYS:HE2	2.17	0.45
1:D:102:ASN:CG	1:D:103:PRO:HD2	2.38	0.45
1:C:198:LEU:HA	1:C:198:LEU:HD23	1.84	0.44
1:D:209:ASP:OD1	1:D:238:LYS:NZ	2.50	0.44
1:D:133:PRO:HG2	1:D:134:GLU:OE2	2.15	0.44
1:C:338:THR:O	1:C:342:LYS:HG3	2.18	0.44
1:B:52:LYS:O	1:B:56:VAL:HG23	2.17	0.44
1:B:284:VAL:O	1:B:288:ILE:HG13	2.16	0.44
1:D:184:ASN:O	1:D:185:ALA:C	2.56	0.44
1:C:135:ARG:NE	4:C:2029:HOH:O	2.49	0.44
1:C:100:LEU:HD13	1:C:322:LYS:HG2	1.99	0.44
1:B:109:ASN:HD22	1:B:150:ALA:HB3	1.81	0.44
1:B:78:ALA:HB1	1:B:178:PRO:CB	2.45	0.44
1:B:102:ASN:CG	1:B:103:PRO:HD2	2.38	0.44
1:C:145:PHE:C	1:C:145:PHE:CD2	2.91	0.44
1:A:145:PHE:CD2	1:A:145:PHE:C	2.91	0.44
1:A:179:LEU:HG	1:A:190:MET:HE1	1.99	0.44
1:C:24:LEU:HD11	1:C:41:ILE:CD1	2.47	0.44
1:B:22:GLU:O	1:B:26:LYS:HG3	2.18	0.44
1:C:63:ARG:HH22	1:C:208:LEU:HB3	1.82	0.44
1:C:58:PHE:HB3	1:C:176:LEU:HD11	1.99	0.44
1:C:71:VAL:HG12	1:C:74:ALA:HB2	2.00	0.44
1:B:24:LEU:HD11	1:B:41:ILE:CD1	2.48	0.43
1:C:239:ARG:NH1	1:C:239:ARG:HG3	2.34	0.43
1:B:332:ASP:OD2	1:B:334:THR:OG1	2.33	0.43
1:D:313:ASP:O	1:D:316:ILE:HG22	2.18	0.43
1:D:110:ARG:HG2	1:D:110:ARG:NH1	2.33	0.43
1:D:95:ALA:HB1	1:D:104:VAL:HG11	2.01	0.43
1:C:137:LYS:CB	4:C:2027:HOH:O	2.61	0.43
1:D:72:PRO:O	1:D:186:LYS:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:VAL:HG23	4:B:2028:HOH:O	2.17	0.43
1:A:115:LYS:HD3	1:A:266:ALA:HB2	2.00	0.43
1:D:179:LEU:HD13	4:D:2013:HOH:O	2.18	0.43
1:D:82:GLY:HA2	1:D:170:ARG:CZ	2.48	0.43
1:B:78:ALA:CB	1:B:188:GLN:HE22	2.31	0.43
1:A:343:SER:CB	4:A:2060:HOH:O	2.67	0.43
1:B:29:ILE:HG22	1:B:156:ALA:HB1	2.01	0.43
1:B:220:PRO:HD3	1:B:228:ILE:CD1	2.49	0.43
1:A:71:VAL:HG12	1:A:74:ALA:HB2	2.01	0.43
1:D:328:SER:CA	1:D:333:VAL:HG22	2.46	0.43
1:C:195:LYS:HB2	1:C:217:TYR:CD2	2.54	0.43
1:A:227:PRO:HG2	1:A:228:ILE:HG23	2.01	0.43
1:B:138:GLU:O	1:B:142:LYS:HG2	2.19	0.43
1:C:102:ASN:CG	1:C:103:PRO:HD2	2.39	0.43
1:A:332:ASP:OD2	1:A:334:THR:OG1	2.33	0.43
1:A:322:LYS:HD2	1:A:325:GLU:HG2	2.01	0.43
1:D:125:LEU:HD13	1:D:270:ALA:HB1	2.00	0.42
1:C:337:LYS:O	1:C:341:VAL:HG23	2.19	0.42
1:D:332:ASP:OD2	1:D:334:THR:OG1	2.34	0.42
1:B:58:PHE:HB3	1:B:176:LEU:HD11	2.01	0.42
1:B:145:PHE:C	1:B:145:PHE:CD2	2.92	0.42
1:C:138:GLU:O	1:C:142:LYS:HG2	2.19	0.42
1:C:187:TYR:HA	1:C:212:LYS:O	2.17	0.42
1:B:115:LYS:HB2	1:B:265:SER:HA	2.02	0.42
1:B:322:LYS:HD2	1:B:325:GLU:HG2	2.02	0.42
1:A:198:LEU:HA	1:A:198:LEU:HD23	1.84	0.42
1:B:90:VAL:HG21	1:B:224:GLU:HG2	2.01	0.42
1:D:282:GLU:HG3	1:D:283:HIS:N	2.35	0.42
1:A:282:GLU:O	1:A:286:GLU:HG3	2.19	0.42
1:C:217:TYR:CZ	1:C:221:GLY:HA2	2.55	0.42
1:D:1:MET:HG2	4:D:2001:HOH:O	2.19	0.42
1:D:58:PHE:HB3	1:D:176:LEU:HD11	2.02	0.42
1:C:87:THR:HG23	1:C:88:VAL:N	2.34	0.42
1:D:337:LYS:O	1:D:341:VAL:HG23	2.20	0.42
1:A:52:LYS:HD2	1:A:207:GLU:OE2	2.19	0.42
1:B:313:ASP:O	1:B:316:ILE:HG22	2.18	0.42
1:A:78:ALA:N	1:A:189:LEU:O	2.52	0.42
1:B:193:PHE:HD2	4:B:2014:HOH:O	2.00	0.42
1:C:220:PRO:HD3	1:C:228:ILE:CD1	2.50	0.42
1:B:83:ASP:O	1:B:84:GLY:C	2.58	0.42
1:A:325:GLU:HG3	1:A:326:ILE:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:239:ARG:NH1	1:D:239:ARG:HG3	2.33	0.42
1:A:26:LYS:HB3	1:A:30:ARG:NH1	2.34	0.42
1:B:54:GLU:O	1:B:58:PHE:HD1	2.03	0.42
1:B:110:ARG:NH1	4:B:2018:HOH:O	2.53	0.42
1:B:78:ALA:N	1:B:189:LEU:O	2.53	0.41
1:D:204:SER:O	1:D:208:LEU:HD23	2.19	0.41
1:A:239:ARG:HG3	1:A:239:ARG:NH1	2.34	0.41
1:C:282:GLU:O	1:C:286:GLU:HG3	2.19	0.41
1:D:280:LYS:HE3	1:D:345:GLY:N	2.15	0.41
1:C:97:LEU:O	1:C:100:LEU:HB2	2.18	0.41
1:B:144:ASN:HD22	1:B:144:ASN:HA	1.65	0.41
1:B:100:LEU:HD13	1:B:322:LYS:HG2	2.02	0.41
1:A:58:PHE:HB3	1:A:176:LEU:HD11	2.02	0.41
1:B:213:ILE:HG23	1:B:236:VAL:HB	2.02	0.41
1:B:190:MET:O	1:B:215:LEU:HA	2.20	0.41
1:A:24:LEU:HD11	1:A:41:ILE:CD1	2.50	0.41
1:D:29:ILE:HG22	1:D:156:ALA:HB1	2.02	0.41
1:D:97:LEU:O	1:D:100:LEU:HB2	2.20	0.41
1:D:26:LYS:HB3	1:D:30:ARG:NH1	2.35	0.41
1:B:282:GLU:HG3	1:B:283:HIS:N	2.36	0.41
1:B:282:GLU:O	1:B:286:GLU:HG3	2.20	0.41
1:A:178:PRO:HG2	1:A:190:MET:HE1	2.02	0.41
1:A:97:LEU:O	1:A:100:LEU:HB2	2.21	0.41
1:A:284:VAL:O	1:A:288:ILE:HG13	2.19	0.41
1:D:213:ILE:HG23	1:D:236:VAL:HB	2.01	0.41
1:D:52:LYS:HE2	1:D:52:LYS:HB3	1.93	0.41
1:D:313:ASP:O	1:D:317:GLU:HG3	2.21	0.41
1:A:100:LEU:HD13	1:A:322:LYS:HG2	2.02	0.41
1:C:313:ASP:O	1:C:317:GLU:HG3	2.20	0.41
1:C:90:VAL:CG2	1:C:224:GLU:HG2	2.51	0.41
1:D:52:LYS:O	1:D:56:VAL:HG23	2.20	0.41
1:B:118:SER:HB3	4:B:2056:HOH:O	2.20	0.41
1:D:154:HIS:HB3	1:D:157:MET:HG2	2.03	0.41
1:B:24:LEU:O	1:B:28:ILE:HG13	2.21	0.41
1:C:26:LYS:HB3	1:C:30:ARG:NH1	2.35	0.41
1:D:196:ASP:O	1:D:197:HIS:C	2.59	0.41
1:C:282:GLU:HG3	1:C:283:HIS:N	2.36	0.41
1:D:154:HIS:HA	4:D:2014:HOH:O	2.20	0.41
1:D:75:ILE:HD12	1:D:299:LEU:HG	2.03	0.41
1:B:51:SER:O	1:B:55:ILE:HG12	2.21	0.41
1:B:233:MET:CG	1:B:244:VAL:HB	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:89:ASN:CB	4:D:2019:HOH:O	2.69	0.41
1:B:313:ASP:O	1:B:317:GLU:HG3	2.20	0.41
1:C:52:LYS:O	1:C:56:VAL:HG23	2.20	0.41
1:B:71:VAL:HG12	1:B:74:ALA:HB2	2.03	0.41
1:C:198:LEU:O	1:C:202:SER:HB3	2.21	0.40
1:A:181:ASN:ND2	1:A:185:ALA:H	2.19	0.40
1:B:191:GLY:HA3	1:B:223:ASP:O	2.20	0.40
1:B:337:LYS:O	1:B:341:VAL:HG23	2.21	0.40
1:C:68:LYS:HG2	4:C:2016:HOH:O	2.21	0.40
1:D:138:GLU:O	1:D:142:LYS:HG2	2.21	0.40
1:C:325:GLU:HG3	1:C:326:ILE:N	2.36	0.40
1:C:29:ILE:HG22	1:C:156:ALA:HB1	2.03	0.40
1:B:328:SER:CA	1:B:333:VAL:HG22	2.46	0.40
1:C:22:GLU:O	1:C:26:LYS:HG3	2.21	0.40
1:A:282:GLU:HG3	1:A:283:HIS:N	2.36	0.40
1:A:90:VAL:CG2	1:A:224:GLU:HG2	2.51	0.40
1:A:337:LYS:O	1:A:341:VAL:HG23	2.21	0.40
1:C:54:GLU:O	1:C:58:PHE:HD1	2.04	0.40
1:B:183:ALA:O	1:B:184:ASN:HB3	2.21	0.40
1:D:282:GLU:O	1:D:286:GLU:HG3	2.20	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	336/345 (97%)	311 (93%)	20 (6%)	5 (2%)	13	28
1	B	336/345 (97%)	311 (93%)	21 (6%)	4 (1%)	16	35
1	C	337/345 (98%)	309 (92%)	23 (7%)	5 (2%)	13	28
1	D	343/345 (99%)	313 (91%)	25 (7%)	5 (2%)	13	28
All	All	1352/1380 (98%)	1244 (92%)	89 (7%)	19 (1%)	14	31

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	SER
1	A	341	VAL
1	B	78	ALA
1	B	341	VAL
1	C	344	SER
1	A	114	GLY
1	A	115	LYS
1	B	84	GLY
1	C	117	GLY
1	D	84	GLY
1	C	113	SER
1	D	115	LYS
1	D	196	ASP
1	D	197	HIS
1	A	330	ASN
1	B	330	ASN
1	C	330	ASN
1	D	330	ASN
1	C	115	LYS

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	285/289 (99%)	274 (96%)	11 (4%)	39	67
1	B	285/289 (99%)	270 (95%)	15 (5%)	28	54
1	C	286/289 (99%)	274 (96%)	12 (4%)	36	64
1	D	288/289 (100%)	276 (96%)	12 (4%)	36	64
All	All	1144/1156 (99%)	1094 (96%)	50 (4%)	35	62

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	VAL

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Mol	Chain	Res	Type
1	A	100	LEU
1	A	125	LEU
1	A	144	ASN
1	A	162	ASN
1	A	184	ASN
1	A	239	ARG
1	A	296	LEU
1	A	306	ARG
1	A	311	TYR
1	A	337	LYS
1	B	87	THR
1	B	90	VAL
1	B	100	LEU
1	B	110	ARG
1	B	125	LEU
1	B	144	ASN
1	B	162	ASN
1	B	193	PHE
1	B	195	LYS
1	B	239	ARG
1	B	296	LEU
1	B	306	ARG
1	B	311	TYR
1	B	337	LYS
1	B	342	LYS
1	C	87	THR
1	C	90	VAL
1	C	100	LEU
1	C	125	LEU
1	C	144	ASN
1	C	162	ASN
1	C	239	ARG
1	C	296	LEU
1	C	301	ARG
1	C	306	ARG
1	C	311	TYR
1	C	337	LYS
1	D	85	LEU
1	D	90	VAL
1	D	100	LEU
1	D	125	LEU
1	D	144	ASN

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Mol	Chain	Res	Type
1	D	162	ASN
1	D	184	ASN
1	D	239	ARG
1	D	296	LEU
1	D	306	ARG
1	D	311	TYR
1	D	337	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN
1	A	151	GLN
1	A	162	ASN
1	A	174	ASN
1	A	181	ASN
1	A	184	ASN
1	A	247	ASN
1	A	283	HIS
1	A	324	ASN
1	B	144	ASN
1	B	151	GLN
1	B	162	ASN
1	B	174	ASN
1	B	181	ASN
1	B	184	ASN
1	B	188	GLN
1	B	247	ASN
1	B	283	HIS
1	B	324	ASN
1	C	109	ASN
1	C	144	ASN
1	C	162	ASN
1	C	174	ASN
1	C	181	ASN
1	C	184	ASN
1	C	247	ASN
1	C	283	HIS
1	C	324	ASN
1	D	144	ASN
1	D	151	GLN
1	D	162	ASN

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Mol	Chain	Res	Type
1	D	174	ASN
1	D	181	ASN
1	D	184	ASN
1	D	188	GLN
1	D	247	ASN
1	D	283	HIS
1	D	324	ASN

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

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
3	POP	A	1346	2	8,8,8	1.98	4 (50%)	13,13,13	1.12	0
3	POP	B	1347	-	8,8,8	1.91	2 (25%)	13,13,13	0.99	1 (7%)
3	POP	C	1348	2	8,8,8	1.54	2 (25%)	13,13,13	0.90	0
3	POP	D	1349	2	8,8,8	2.05	3 (37%)	13,13,13	1.01	1 (7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	POP	A	1346	2	-	0/6/6/6	0/0/0/0
3	POP	B	1347	-	-	0/6/6/6	0/0/0/0
3	POP	C	1348	2	-	0/6/6/6	0/0/0/0
3	POP	D	1349	2	-	0/6/6/6	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1347	POP	P2-O	-2.20	1.56	1.61
3	A	1346	POP	P1-O3	2.03	1.57	1.50
3	C	1348	POP	P2-O6	2.27	1.58	1.50
3	A	1346	POP	P2-O6	2.51	1.58	1.50
3	A	1346	POP	P1-O	2.59	1.67	1.61
3	C	1348	POP	P1-O3	2.66	1.59	1.50
3	D	1349	POP	P2-O6	2.89	1.60	1.50
3	D	1349	POP	P1-O1	2.95	1.59	1.50
3	A	1346	POP	P1-O1	3.23	1.60	1.50
3	D	1349	POP	P1-O3	3.65	1.62	1.50
3	B	1347	POP	P1-O3	3.80	1.63	1.50

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	1349	POP	P2-O-P1	-2.74	125.03	132.73
3	B	1347	POP	P2-O-P1	-2.66	125.25	132.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1347	POP	2	0
3	C	1348	POP	1	0
3	D	1349	POP	4	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	340/345 (98%)	0.88	53 (15%)	3 2	34, 67, 120, 146	0
1	B	340/345 (98%)	0.73	40 (11%)	6 4	33, 68, 122, 174	0
1	C	341/345 (98%)	0.71	35 (10%)	9 6	37, 66, 115, 175	0
1	D	345/345 (100%)	1.08	65 (18%)	2 1	36, 76, 131, 177	0
All	All	1366/1380 (98%)	0.85	193 (14%)	4 2	33, 69, 125, 177	0

All (193) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	SER	8.1
1	D	82	GLY	7.5
1	B	83	ASP	7.2
1	D	345	GLY	7.0
1	D	83	ASP	7.0
1	D	85	LEU	6.8
1	C	345	GLY	6.7
1	A	113	SER	6.3
1	B	112	VAL	5.8
1	C	112	VAL	5.7
1	B	1	MET	5.3
1	D	112	VAL	5.2
1	B	141	ASN	5.0
1	A	112	VAL	5.0
1	D	142	LYS	4.8
1	D	239	ARG	4.8
1	D	252	GLY	4.8
1	C	302	VAL	4.8
1	A	135	ARG	4.5
1	D	113	SER	4.4
1	C	314	HIS	4.4

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Mol	Chain	Res	Type	RSRZ
1	D	290	ILE	4.3
1	D	280	LYS	4.3
1	C	115	LYS	4.2
1	D	257	PRO	4.2
1	B	113	SER	4.1
1	C	291	ASN	4.0
1	D	306	ARG	4.0
1	D	84	GLY	4.0
1	A	157	MET	4.0
1	B	114	GLY	4.0
1	B	239	ARG	3.9
1	D	282	GLU	3.8
1	D	80	THR	3.8
1	A	250	ASP	3.7
1	C	85	LEU	3.7
1	D	114	GLY	3.7
1	A	12	ASN	3.7
1	B	111	ALA	3.6
1	C	239	ARG	3.6
1	A	170	ARG	3.5
1	B	156	ALA	3.5
1	B	280	LYS	3.5
1	A	152	TYR	3.5
1	B	328	SER	3.5
1	A	1	MET	3.5
1	D	233	MET	3.4
1	D	215	LEU	3.3
1	B	158	LYS	3.3
1	A	240	GLY	3.3
1	D	243	GLU	3.3
1	B	115	LYS	3.3
1	C	113	SER	3.2
1	D	193	PHE	3.2
1	A	15	ASP	3.2
1	D	307	GLU	3.2
1	B	157	MET	3.2
1	D	245	LYS	3.2
1	B	116	SER	3.2
1	D	40	ALA	3.2
1	B	138	GLU	3.2
1	A	239	ARG	3.2
1	D	206	TYR	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	225	VAL	3.1
1	D	275	ARG	3.1
1	C	282	GLU	3.1
1	A	343	SER	3.1
1	A	196	ASP	3.1
1	C	344	SER	3.1
1	B	341	VAL	3.1
1	A	243	GLU	3.1
1	C	303	GLY	3.0
1	A	233	MET	3.0
1	D	115	LYS	3.0
1	C	290	ILE	3.0
1	A	138	GLU	2.9
1	A	254	SER	2.9
1	A	234	LYS	2.9
1	C	246	LEU	2.8
1	D	228	ILE	2.8
1	B	196	ASP	2.8
1	B	334	THR	2.8
1	A	232	PHE	2.8
1	D	291	ASN	2.8
1	D	238	LYS	2.8
1	D	1	MET	2.8
1	B	232	PHE	2.8
1	B	249	THR	2.8
1	C	211	ASN	2.8
1	B	110	ARG	2.8
1	A	115	LYS	2.7
1	D	279	GLY	2.7
1	B	14	SER	2.7
1	A	217	TYR	2.7
1	D	227	PRO	2.7
1	B	245	LYS	2.7
1	A	249	THR	2.7
1	C	269	SER	2.7
1	D	344	SER	2.6
1	D	41	ILE	2.6
1	D	283	HIS	2.6
1	D	317	GLU	2.6
1	B	159	ASN	2.6
1	D	303	GLY	2.6
1	C	210	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	2	ASN	2.6
1	B	195	LYS	2.6
1	D	78	ALA	2.6
1	A	303	GLY	2.5
1	A	206	TYR	2.5
1	A	252	GLY	2.5
1	A	300	ASP	2.5
1	C	40	ALA	2.5
1	C	84	GLY	2.5
1	D	287	PHE	2.5
1	D	2	ASN	2.5
1	C	111	ALA	2.5
1	C	41	ILE	2.5
1	D	273	ILE	2.5
1	D	271	ILE	2.5
1	D	292	THR	2.5
1	C	101	VAL	2.4
1	D	172	ILE	2.4
1	C	334	THR	2.4
1	A	40	ALA	2.4
1	D	294	VAL	2.4
1	A	245	LYS	2.4
1	A	221	GLY	2.4
1	C	213	ILE	2.4
1	A	114	GLY	2.4
1	A	70	ASP	2.4
1	A	199	ASP	2.4
1	C	170	ARG	2.3
1	C	116	SER	2.3
1	D	314	HIS	2.3
1	D	285	ALA	2.3
1	C	273	ILE	2.3
1	B	72	PRO	2.3
1	C	78	ALA	2.3
1	C	270	ALA	2.3
1	B	109	ASN	2.3
1	A	155	PRO	2.3
1	A	202	SER	2.3
1	B	266	ALA	2.3
1	B	19	ASN	2.3
1	A	134	GLU	2.2
1	D	190	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	231	THR	2.2
1	B	151	GLN	2.2
1	B	134	GLU	2.2
1	B	155	PRO	2.2
1	C	280	LYS	2.2
1	C	168	GLY	2.2
1	D	309	TYR	2.2
1	A	2	ASN	2.2
1	C	306	ARG	2.2
1	A	27	ALA	2.2
1	A	156	ALA	2.2
1	D	214	ILE	2.2
1	D	254	SER	2.2
1	D	251	PHE	2.2
1	A	230	ASN	2.2
1	D	260	LYS	2.1
1	D	264	ASN	2.1
1	D	310	GLU	2.1
1	A	13	LYS	2.1
1	A	159	ASN	2.1
1	D	138	GLU	2.1
1	A	205	ALA	2.1
1	D	42	LEU	2.1
1	B	135	ARG	2.1
1	A	187	TYR	2.1
1	A	38	VAL	2.1
1	A	341	VAL	2.1
1	D	216	VAL	2.1
1	B	238	LYS	2.1
1	D	334	THR	2.1
1	D	79	GLY	2.1
1	A	37	LEU	2.1
1	A	42	LEU	2.1
1	A	195	LYS	2.0
1	C	193	PHE	2.0
1	C	238	LYS	2.0
1	D	288	ILE	2.0
1	D	116	SER	2.0
1	C	147	PHE	2.0
1	A	78	ALA	2.0
1	A	109	ASN	2.0
1	B	30	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	146	VAL	2.0
1	D	96	ILE	2.0
1	B	154	HIS	2.0
1	B	265	SER	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	POP	A	1346	9/9	0.91	0.26	0.90	21,21,21,21	0
3	POP	B	1347	9/9	0.95	0.23	0.44	21,21,21,76	0
3	POP	C	1348	9/9	0.97	0.21	-0.11	21,21,21,21	0
2	MG	A	1344	1/1	0.57	0.13	-1.35	42,42,42,42	0
3	POP	D	1349	9/9	0.96	0.14	-1.72	21,21,21,21	0
2	MG	D	1350	1/1	0.98	0.05	-3.17	59,59,59,59	0
2	MG	C	1346	1/1	0.88	0.13	-	47,47,47,47	0

6.5 Other polymers (i)

There are no such residues in this entry.