Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 09:46 AM BST

PDB ID	:	4GJT
Title	:	complex structure of nectin-4 bound to MV-H
Authors	:	Zhang, X.; Lu, G.; Qi, J.; Li, Y.; He, Y.; Xu, X.; Shi, J.; Zhang, C.; Yan, J.; Gao, G.F.
Deposited on	:	2012-08-10
Resolution	:	3.10 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

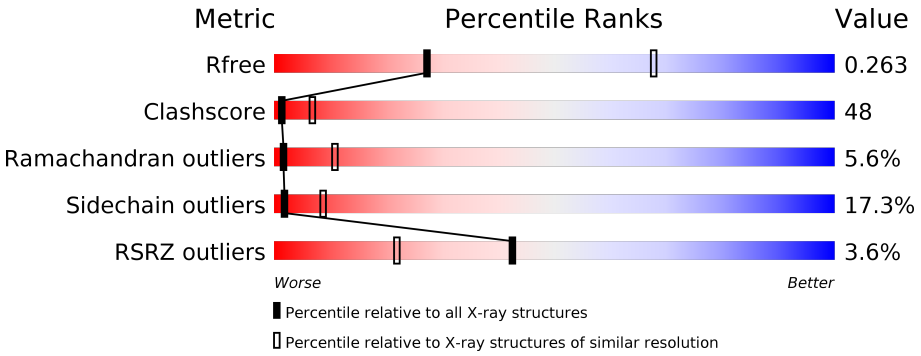
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

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 3.10 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	473	2% 35% 39% 11% • 13%
2	B	123	8% 40% 45% 7% • 7%
2	C	123	3% 49% 36% 7% 9%

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
3	NAG	A	801	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5020 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 Hemagglutinin glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3205	2051	540	591	23			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	ALA	-	expression tag	UNP Q786F2
A	146	ASP	-	expression tag	UNP Q786F2
A	147	GLY	-	expression tag	UNP Q786F2
A	148	ILE	-	expression tag	UNP Q786F2
A	149	GLN	-	expression tag	UNP Q786F2
A	150	HIS	-	expression tag	UNP Q786F2
A	151	HIS	-	expression tag	UNP Q786F2
A	152	HIS	-	expression tag	UNP Q786F2
A	153	HIS	-	expression tag	UNP Q786F2
A	154	HIS	-	expression tag	UNP Q786F2
A	155	HIS	-	expression tag	UNP Q786F2

- Molecule 2 is a protein called Poliovirus receptor-related protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	114	Total	C	N	O	S	0	0	0
			870	542	157	169	2			
2	C	112	Total	C	N	O	S	0	0	0
			852	531	152	167	2			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	expression tag	UNP Q96NY8
B	116	GLU	-	expression tag	UNP Q96NY8
B	117	HIS	-	expression tag	UNP Q96NY8

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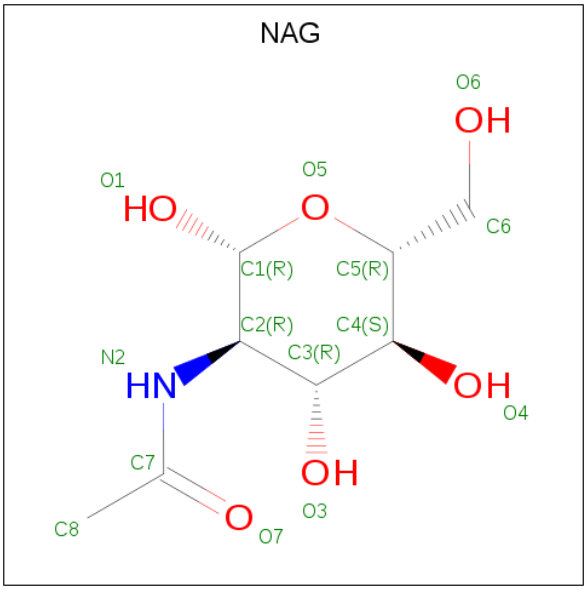
Chain	Residue	Modelled	Actual	Comment	Reference
B	118	HIS	-	expression tag	UNP Q96NY8
B	119	HIS	-	expression tag	UNP Q96NY8
B	120	HIS	-	expression tag	UNP Q96NY8
B	121	HIS	-	expression tag	UNP Q96NY8
B	122	HIS	-	expression tag	UNP Q96NY8
C	0	MET	-	expression tag	UNP Q96NY8
C	116	GLU	-	expression tag	UNP Q96NY8
C	117	HIS	-	expression tag	UNP Q96NY8
C	118	HIS	-	expression tag	UNP Q96NY8
C	119	HIS	-	expression tag	UNP Q96NY8
C	120	HIS	-	expression tag	UNP Q96NY8
C	121	HIS	-	expression tag	UNP Q96NY8
C	122	HIS	-	expression tag	UNP Q96NY8

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	63	Total	O	0	0
			63	63		

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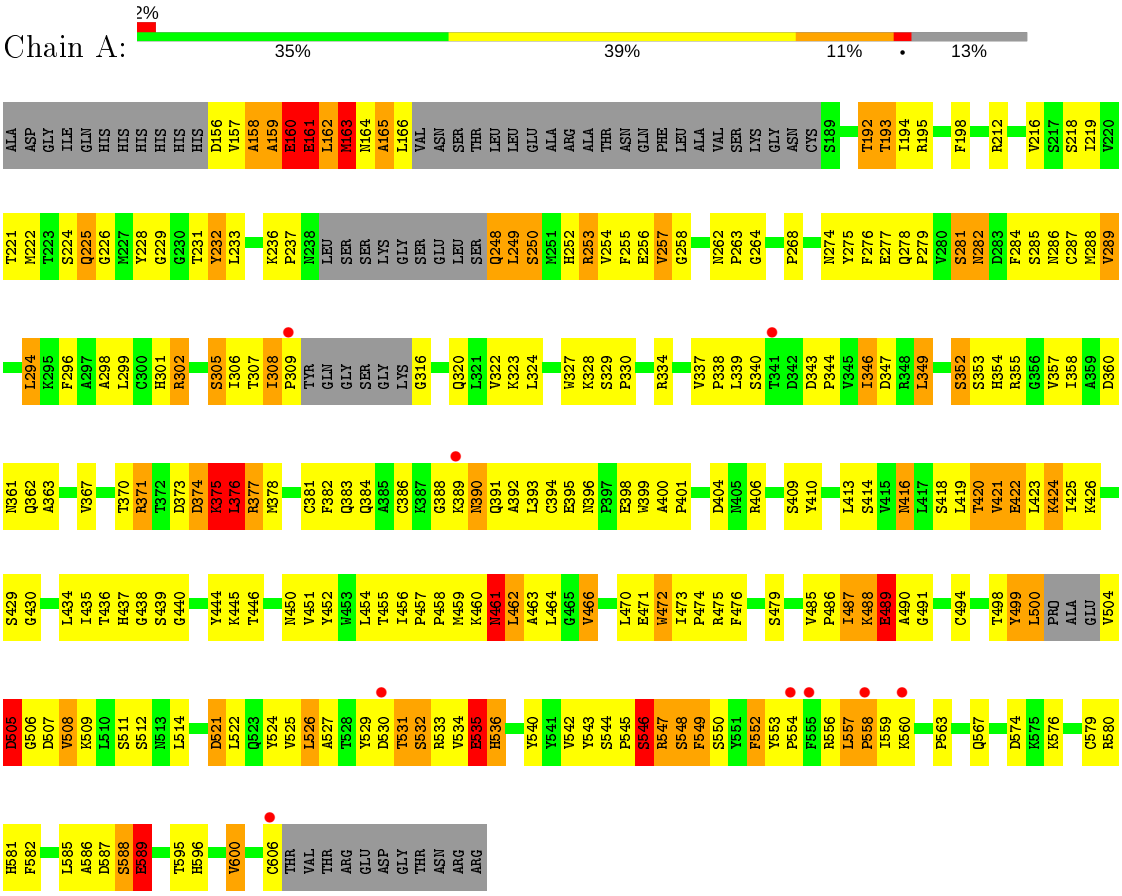
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	6	Total O 6 6	0	0
4	C	10	Total O 10 10	0	0

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.

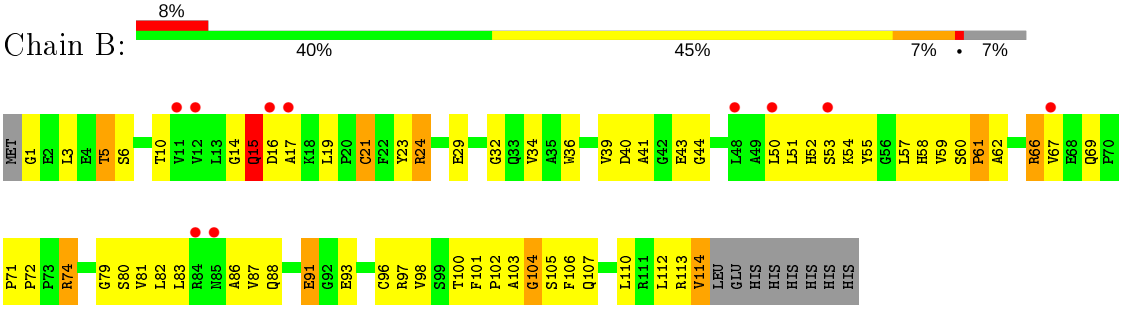
- Molecule 1: Hemagglutinin glycoprotein

Chain A:



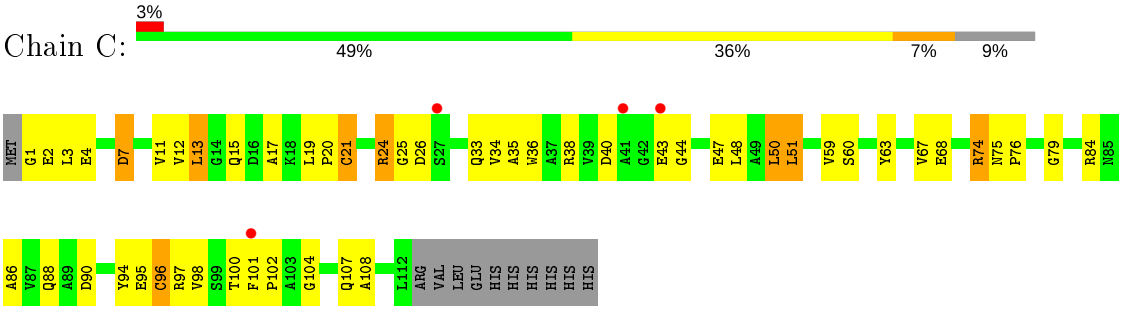
- Molecule 2: Poliovirus receptor-related protein 4

Chain B:



• Molecule 2: Poliovirus receptor-related protein 4

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	82.35Å 171.78Å 117.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.13 – 3.10 37.13 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.13-3.10) 99.9 (37.13-3.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.21 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.260 , 0.278 0.254 , 0.263	Depositor DCC
R_{free} test set	774 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	90.8	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 59.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5020	wwPDB-VP
Average B, all atoms (Å^2)	96.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NAG

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	0.35	0/3285	0.57	4/4465 (0.1%)
2	B	0.28	0/888	0.49	0/1207
2	C	0.39	0/870	0.50	0/1183
All	All	0.35	0/5043	0.55	4/6855 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	165	ALA	CB-CA-C	-7.93	98.20	110.10
1	A	586	ALA	CB-CA-C	7.37	121.16	110.10
1	A	165	ALA	N-CA-C	6.81	129.38	111.00
1	A	505	ASP	CB-CA-C	-5.23	99.94	110.40

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	3205	0	3153	370	2
2	B	870	0	844	66	2
2	C	852	0	822	44	0
3	A	14	0	13	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	63	0	0	62	0
4	B	6	0	0	9	0
4	C	10	0	0	9	0
All	All	5020	0	4832	472	2

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 48.

All (472) 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:A:489:GLU:HB2	4:A:958:HOH:O	1.24	1.33
1:A:464:LEU:CD2	2:B:100:THR:O	1.77	1.30
1:A:588:SER:O	1:A:589:GLU:HG3	1.36	1.22
1:A:376:LEU:N	1:A:376:LEU:HD23	1.55	1.19
2:B:54:LYS:HG3	2:B:55:TYR:CD1	1.83	1.12
1:A:163:MET:HG3	1:A:164:ASN:N	1.54	1.11
1:A:420:THR:HA	1:A:421:VAL:HG23	1.14	1.11
1:A:156:ASP:HA	4:A:939:HOH:O	1.47	1.11
1:A:376:LEU:H	1:A:376:LEU:CD2	1.58	1.10
1:A:588:SER:O	1:A:589:GLU:CG	1.99	1.10
1:A:160:GLU:HG3	1:A:161:GLU:N	1.41	1.09
1:A:390:ASN:HA	1:A:499:TYR:CD2	1.86	1.08
1:A:588:SER:C	1:A:589:GLU:HG3	1.69	1.08
1:A:499:TYR:OH	2:B:54:LYS:HE2	1.54	1.07
1:A:323:LYS:NZ	4:A:943:HOH:O	1.86	1.07
2:B:54:LYS:HG3	2:B:55:TYR:HD1	1.10	1.06
1:A:163:MET:O	1:A:166:LEU:HD23	1.53	1.06
1:A:461:ASN:N	4:A:907:HOH:O	1.86	1.06
1:A:489:GLU:N	4:A:958:HOH:O	1.87	1.05
1:A:464:LEU:HD21	2:B:100:THR:O	0.88	1.04
1:A:534:VAL:O	1:A:534:VAL:HG12	1.57	1.04
1:A:437:HIS:HB3	1:A:459:MET:HE2	1.36	1.04
1:A:360:ASP:CG	3:A:801:NAG:O6	1.98	1.02
1:A:504:VAL:HG13	1:A:504:VAL:O	1.58	1.01
2:B:54:LYS:HE3	2:B:55:TYR:HE1	1.25	1.01
1:A:420:THR:HA	1:A:421:VAL:CG2	1.88	1.01
1:A:532:SER:HA	4:A:921:HOH:O	1.61	0.99
1:A:500:LEU:HD23	2:B:32:GLY:HA2	1.43	0.99
1:A:376:LEU:H	1:A:376:LEU:HD23	0.81	0.97
1:A:437:HIS:HB3	1:A:459:MET:CE	1.94	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ALA:O	1:A:162:LEU:HB2	1.66	0.95
1:A:262:ASN:OD1	4:A:928:HOH:O	1.83	0.95
1:A:438:GLY:C	4:A:922:HOH:O	2.05	0.94
1:A:509:LYS:NZ	1:A:531:THR:HG21	1.83	0.94
1:A:587:ASP:OD1	4:A:929:HOH:O	1.84	0.93
1:A:281:SER:OG	4:A:932:HOH:O	1.85	0.93
2:B:74:ARG:HB2	2:B:74:ARG:HH11	1.34	0.92
1:A:160:GLU:CG	1:A:161:GLU:N	2.30	0.92
2:C:74:ARG:HB3	4:C:201:HOH:O	1.71	0.91
1:A:160:GLU:HG3	1:A:161:GLU:H	1.05	0.91
1:A:355:ARG:NH2	1:A:567:GLN:HG3	1.85	0.91
1:A:340:SER:HB2	1:A:424:LYS:HG2	1.52	0.90
1:A:375:LYS:N	1:A:375:LYS:CD	2.30	0.90
1:A:418:SER:O	1:A:419:LEU:HB2	1.71	0.90
1:A:472:TRP:O	1:A:475:ARG:N	2.04	0.90
1:A:399:TRP:CB	4:A:924:HOH:O	2.20	0.90
1:A:488:LYS:O	1:A:490:ALA:N	2.04	0.89
1:A:163:MET:CG	1:A:164:ASN:N	2.35	0.89
1:A:461:ASN:O	1:A:462:LEU:HG	1.72	0.89
1:A:198:PHE:HZ	1:A:554:PRO:HD2	1.38	0.89
2:C:4:GLU:OE1	4:C:207:HOH:O	1.91	0.89
1:A:163:MET:HG3	1:A:164:ASN:H	1.35	0.88
2:B:44:GLY:HA2	4:B:203:HOH:O	1.73	0.88
1:A:458:PRO:O	4:A:910:HOH:O	1.91	0.88
2:C:97:ARG:HG2	2:C:107:GLN:HG3	1.56	0.88
1:A:524:TYR:CE1	1:A:543:TYR:CD1	2.62	0.88
1:A:375:LYS:HD3	1:A:375:LYS:H	1.37	0.87
1:A:160:GLU:CG	1:A:161:GLU:H	1.86	0.87
1:A:162:LEU:C	4:A:926:HOH:O	2.13	0.85
1:A:418:SER:N	4:A:938:HOH:O	2.07	0.85
2:C:21:CYS:HB2	2:C:36:TRP:HZ2	1.41	0.85
1:A:438:GLY:O	4:A:922:HOH:O	1.94	0.85
2:B:74:ARG:NH1	2:B:74:ARG:HB2	1.91	0.84
3:A:801:NAG:C3	4:A:902:HOH:O	2.26	0.84
1:A:159:ALA:O	1:A:162:LEU:CB	2.25	0.83
3:A:801:NAG:O3	4:A:902:HOH:O	1.95	0.83
1:A:375:LYS:H	1:A:375:LYS:CD	1.88	0.81
1:A:390:ASN:CA	1:A:499:TYR:CD2	2.62	0.81
1:A:288:MET:HE3	1:A:299:LEU:HB3	1.62	0.81
3:A:801:NAG:H4	4:A:902:HOH:O	1.80	0.81
2:B:54:LYS:HE3	2:B:55:TYR:CE1	2.12	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:MET:CG	1:A:164:ASN:H	1.93	0.80
1:A:192:THR:O	1:A:193:THR:CG2	2.30	0.80
1:A:534:VAL:O	1:A:534:VAL:CG1	2.30	0.80
1:A:163:MET:N	4:A:926:HOH:O	2.15	0.80
1:A:504:VAL:CG1	1:A:504:VAL:O	2.30	0.79
1:A:160:GLU:O	1:A:162:LEU:N	2.16	0.79
1:A:157:VAL:O	1:A:159:ALA:N	2.16	0.79
1:A:549:PHE:C	1:A:549:PHE:CD2	2.56	0.79
1:A:488:LYS:C	1:A:490:ALA:H	1.87	0.78
2:C:59:VAL:O	4:C:209:HOH:O	2.01	0.78
1:A:163:MET:CA	4:A:926:HOH:O	2.30	0.78
1:A:355:ARG:HH11	1:A:355:ARG:HG3	1.49	0.78
1:A:511:SER:OG	4:A:910:HOH:O	1.99	0.78
1:A:461:ASN:HB2	1:A:508:VAL:O	1.85	0.77
1:A:509:LYS:HZ2	1:A:531:THR:HG21	1.47	0.77
1:A:437:HIS:CB	1:A:459:MET:HE2	2.14	0.77
2:B:34:VAL:HG13	2:B:98:VAL:HG22	1.67	0.77
1:A:157:VAL:HG12	1:A:158:ALA:N	2.00	0.77
1:A:450:ASN:CG	1:A:472:TRP:HE1	1.88	0.77
1:A:285:SER:HB3	4:A:912:HOH:O	1.85	0.76
1:A:557:LEU:CB	1:A:558:PRO:HD2	2.15	0.76
1:A:461:ASN:O	1:A:462:LEU:CG	2.34	0.76
1:A:552:PHE:O	1:A:554:PRO:HD3	1.85	0.76
2:C:21:CYS:HB2	2:C:36:TRP:CZ2	2.21	0.76
1:A:420:THR:CA	1:A:421:VAL:HG23	2.06	0.75
1:A:250:SER:HB3	1:A:252:HIS:CE1	2.21	0.75
1:A:437:HIS:HD2	1:A:439:SER:H	1.32	0.75
2:C:4:GLU:OE2	2:C:24:ARG:HD3	1.86	0.75
1:A:450:ASN:HA	1:A:472:TRP:CZ2	2.21	0.75
1:A:375:LYS:N	1:A:375:LYS:HD3	1.96	0.75
1:A:221:THR:HG22	1:A:567:GLN:HE22	1.52	0.74
1:A:228:TYR:CE1	1:A:258:GLY:HA3	2.23	0.74
1:A:536:HIS:O	1:A:556:ARG:HG2	1.87	0.74
2:B:112:LEU:O	4:B:202:HOH:O	2.04	0.74
2:B:105:SER:OG	4:B:206:HOH:O	2.04	0.74
1:A:500:LEU:CD2	2:B:32:GLY:HA2	2.18	0.73
1:A:163:MET:HA	4:A:926:HOH:O	1.86	0.73
2:B:112:LEU:HB3	4:B:202:HOH:O	1.87	0.73
1:A:374:ASP:HB2	1:A:375:LYS:HD3	1.70	0.73
1:A:198:PHE:CZ	1:A:554:PRO:HD2	2.22	0.72
1:A:560:LYS:HB3	4:A:929:HOH:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:SER:CA	4:A:938:HOH:O	2.37	0.72
1:A:218:SER:HB2	1:A:233:LEU:HD23	1.70	0.72
1:A:237:PRO:HA	1:A:249:LEU:CD1	2.20	0.72
1:A:452:TYR:CD2	1:A:472:TRP:CH2	2.78	0.72
1:A:499:TYR:OH	2:B:54:LYS:CE	2.35	0.72
1:A:452:TYR:CD2	1:A:472:TRP:HH2	2.07	0.71
1:A:248:GLN:O	1:A:249:LEU:HD12	1.91	0.71
1:A:532:SER:CA	4:A:921:HOH:O	2.25	0.71
1:A:459:MET:O	4:A:907:HOH:O	2.07	0.71
1:A:399:TRP:HB3	4:A:924:HOH:O	1.83	0.71
1:A:390:ASN:H	1:A:499:TYR:HD2	1.38	0.71
3:A:801:NAG:H2	4:A:902:HOH:O	1.90	0.71
1:A:166:LEU:HD12	1:A:166:LEU:O	1.90	0.71
1:A:535:GLU:O	1:A:536:HIS:O	2.08	0.71
1:A:489:GLU:CB	4:A:958:HOH:O	1.98	0.70
1:A:355:ARG:HH21	1:A:567:GLN:HG3	1.52	0.70
1:A:500:LEU:HD23	2:B:32:GLY:CA	2.21	0.70
2:B:91:GLU:HG3	2:B:114:VAL:HG23	1.73	0.70
1:A:529:TYR:CD2	1:A:563:PRO:HG3	2.27	0.70
1:A:162:LEU:O	4:A:926:HOH:O	2.10	0.70
2:B:34:VAL:HG21	2:B:79:GLY:HA3	1.72	0.70
1:A:156:ASP:CG	1:A:157:VAL:H	1.94	0.69
1:A:509:LYS:HZ1	1:A:531:THR:HG21	1.57	0.69
1:A:595:THR:O	4:A:949:HOH:O	2.10	0.69
2:C:67:VAL:O	4:C:210:HOH:O	2.10	0.69
1:A:488:LYS:C	1:A:490:ALA:N	2.45	0.68
2:C:34:VAL:HG22	2:C:98:VAL:HG22	1.73	0.68
1:A:531:THR:O	4:A:921:HOH:O	2.12	0.67
1:A:450:ASN:OD1	1:A:472:TRP:NE1	2.20	0.67
1:A:416:ASN:OD1	3:A:801:NAG:H2	1.94	0.67
1:A:285:SER:CB	4:A:912:HOH:O	2.40	0.67
1:A:418:SER:OG	4:A:938:HOH:O	2.13	0.67
1:A:544:SER:OG	1:A:545:PRO:HD2	1.95	0.67
2:C:3:LEU:HD13	2:C:108:ALA:HB2	1.75	0.67
2:C:7:ASP:O	4:C:205:HOH:O	2.11	0.67
1:A:192:THR:O	1:A:193:THR:HG23	1.94	0.66
2:B:50:LEU:HB3	2:B:58:HIS:HB3	1.76	0.66
1:A:410:TYR:CE2	1:A:435:ILE:HD11	2.30	0.66
1:A:588:SER:O	1:A:589:GLU:HG2	1.95	0.66
1:A:375:LYS:HB2	1:A:376:LEU:HD23	1.78	0.66
1:A:450:ASN:CG	1:A:472:TRP:NE1	2.48	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:ASN:N	1:A:499:TYR:CD2	2.64	0.66
1:A:524:TYR:CE1	1:A:543:TYR:HD1	2.12	0.66
3:A:801:NAG:O3	4:A:931:HOH:O	2.13	0.66
1:A:504:VAL:N	4:A:915:HOH:O	2.28	0.66
1:A:237:PRO:HA	1:A:249:LEU:HD11	1.78	0.65
1:A:535:GLU:O	1:A:536:HIS:C	2.33	0.65
1:A:160:GLU:C	1:A:162:LEU:N	2.50	0.65
1:A:399:TRP:N	4:A:924:HOH:O	2.19	0.65
1:A:461:ASN:O	1:A:461:ASN:ND2	2.30	0.65
1:A:487:ILE:HG22	1:A:487:ILE:O	1.96	0.65
1:A:505:ASP:N	1:A:505:ASP:OD2	2.30	0.65
1:A:156:ASP:OD1	1:A:157:VAL:N	2.30	0.65
1:A:157:VAL:O	1:A:158:ALA:C	2.34	0.65
2:B:44:GLY:CA	4:B:203:HOH:O	2.38	0.64
1:A:192:THR:C	1:A:193:THR:HG22	2.18	0.64
1:A:420:THR:HA	1:A:421:VAL:CB	2.22	0.64
1:A:316:GLY:O	2:C:24:ARG:NH2	2.30	0.64
1:A:505:ASP:O	4:A:960:HOH:O	2.15	0.64
1:A:192:THR:O	1:A:193:THR:HG22	1.97	0.64
1:A:546:SER:O	4:A:906:HOH:O	2.15	0.64
1:A:381:CYS:HA	1:A:490:ALA:HB1	1.80	0.63
1:A:420:THR:O	4:A:946:HOH:O	2.15	0.63
1:A:361:ASN:ND2	1:A:418:SER:HB3	2.14	0.63
1:A:248:GLN:O	1:A:249:LEU:CD1	2.47	0.63
1:A:472:TRP:O	1:A:475:ARG:O	2.16	0.63
1:A:299:LEU:HD11	1:A:425:ILE:HD11	1.81	0.62
1:A:461:ASN:C	1:A:462:LEU:HG	2.19	0.62
2:B:5:THR:CB	4:B:201:HOH:O	2.40	0.62
1:A:354:HIS:CE1	1:A:367:VAL:HG12	2.34	0.62
2:B:54:LYS:HG3	2:B:55:TYR:CE1	2.33	0.62
1:A:549:PHE:C	1:A:549:PHE:HD2	2.02	0.62
1:A:390:ASN:ND2	1:A:436:THR:HG22	2.15	0.62
2:C:3:LEU:HD21	2:C:98:VAL:HG21	1.81	0.61
1:A:302:ARG:NH1	4:A:944:HOH:O	2.32	0.61
2:C:40:ASP:OD1	2:C:44:GLY:HA2	2.00	0.61
1:A:250:SER:HB3	1:A:252:HIS:HE1	1.62	0.61
1:A:507:ASP:O	1:A:531:THR:HG23	2.01	0.61
1:A:373:ASP:HB2	1:A:376:LEU:HG	1.83	0.61
1:A:163:MET:O	1:A:166:LEU:CD2	2.40	0.60
1:A:360:ASP:CB	3:A:801:NAG:O6	2.48	0.60
1:A:157:VAL:CG1	1:A:158:ALA:N	2.64	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ASP:OD1	3:A:801:NAG:O6	2.19	0.60
1:A:375:LYS:N	1:A:375:LYS:HD2	2.13	0.60
1:A:384:GLN:OE1	1:A:489:GLU:HB3	2.01	0.60
2:B:44:GLY:N	4:B:203:HOH:O	2.34	0.60
1:A:587:ASP:HA	4:A:929:HOH:O	2.01	0.60
1:A:375:LYS:O	1:A:376:LEU:C	2.40	0.60
1:A:281:SER:N	4:A:932:HOH:O	2.33	0.59
1:A:557:LEU:CB	1:A:558:PRO:CD	2.80	0.59
1:A:461:ASN:O	1:A:462:LEU:CD2	2.50	0.59
2:B:3:LEU:HD11	2:B:98:VAL:HG23	1.84	0.59
1:A:374:ASP:O	1:A:375:LYS:O	2.20	0.59
2:B:67:VAL:HA	2:B:82:LEU:O	2.03	0.59
2:B:112:LEU:HD12	2:B:113:ARG:H	1.67	0.59
1:A:221:THR:H	1:A:567:GLN:NE2	2.00	0.58
1:A:419:LEU:O	1:A:421:VAL:HG23	2.03	0.58
1:A:526:LEU:HG	1:A:527:ALA:N	2.18	0.58
1:A:398:GLU:HG3	1:A:398:GLU:O	2.03	0.58
1:A:355:ARG:HH21	1:A:567:GLN:HE21	1.51	0.58
1:A:444:TYR:OH	4:A:916:HOH:O	2.00	0.58
2:B:60:SER:HB3	2:B:61:PRO:HD2	1.86	0.57
2:B:21:CYS:HB2	2:B:36:TRP:CZ2	2.39	0.57
1:A:499:TYR:CZ	2:B:54:LYS:HE2	2.39	0.57
1:A:532:SER:C	4:A:921:HOH:O	2.41	0.57
1:A:221:THR:CG2	1:A:567:GLN:HE22	2.18	0.57
1:A:289:VAL:CG2	1:A:296:PHE:HE1	2.18	0.57
1:A:371:ARG:HH11	1:A:429:SER:HB3	1.69	0.57
1:A:399:TRP:CA	4:A:924:HOH:O	2.52	0.57
1:A:409:SER:HB2	1:A:430:GLY:O	2.04	0.57
1:A:521:ASP:HB2	1:A:545:PRO:HG2	1.84	0.57
1:A:371:ARG:HD3	1:A:429:SER:OG	2.04	0.56
1:A:390:ASN:HA	1:A:499:TYR:CE2	2.38	0.56
1:A:236:LYS:HD2	1:A:252:HIS:CE1	2.40	0.56
1:A:257:VAL:HG13	1:A:274:ASN:HB3	1.85	0.56
1:A:547:ARG:HD3	1:A:547:ARG:C	2.25	0.56
1:A:253:ARG:HG3	1:A:253:ARG:HH11	1.70	0.56
3:A:801:NAG:C2	4:A:902:HOH:O	2.47	0.56
1:A:534:VAL:O	1:A:535:GLU:HB2	2.06	0.56
2:C:38:ARG:HG3	2:C:48:LEU:HD11	1.87	0.56
1:A:192:THR:C	1:A:193:THR:CG2	2.74	0.56
2:C:33:GLN:CD	2:C:50:LEU:HD11	2.26	0.56
1:A:525:VAL:HG22	1:A:526:LEU:N	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526:LEU:HD21	2:B:102:PRO:HB2	1.87	0.56
1:A:391:GLN:O	1:A:392:ALA:HB3	2.06	0.56
2:C:63:TYR:HB3	2:C:67:VAL:HG23	1.88	0.55
1:A:524:TYR:HE1	1:A:543:TYR:CD1	2.21	0.55
1:A:231:THR:HA	1:A:254:VAL:O	2.06	0.55
1:A:355:ARG:HH21	1:A:567:GLN:CG	2.19	0.55
1:A:543:TYR:CE1	2:B:103:ALA:HB3	2.42	0.55
1:A:461:ASN:O	1:A:462:LEU:HD23	2.07	0.55
1:A:288:MET:O	1:A:298:ALA:HA	2.07	0.54
1:A:534:VAL:O	1:A:535:GLU:CB	2.55	0.54
1:A:529:TYR:OH	1:A:536:HIS:CD2	2.60	0.54
1:A:390:ASN:ND2	1:A:436:THR:CG2	2.71	0.54
2:C:48:LEU:HD21	2:C:94:TYR:HE1	1.72	0.54
1:A:375:LYS:O	1:A:377:ARG:N	2.40	0.54
1:A:446:THR:OG1	1:A:451:VAL:HG13	2.07	0.54
1:A:544:SER:OG	1:A:545:PRO:CD	2.56	0.54
1:A:500:LEU:O	4:A:954:HOH:O	2.17	0.54
1:A:461:ASN:OD1	1:A:506:GLY:HA3	2.07	0.54
1:A:529:TYR:OH	1:A:536:HIS:HD2	1.90	0.54
1:A:547:ARG:HD3	1:A:548:SER:N	2.23	0.53
1:A:221:THR:H	1:A:567:GLN:HE22	1.56	0.53
1:A:255:PHE:CZ	1:A:322:VAL:HG21	2.44	0.53
1:A:445:LYS:HG2	1:A:452:TYR:HE1	1.72	0.53
4:A:917:HOH:O	2:B:54:LYS:HD3	2.08	0.53
2:B:100:THR:OG1	2:B:104:GLY:HA3	2.08	0.53
1:A:327:TRP:O	1:A:328:LYS:C	2.47	0.53
2:B:1:GLY:HA3	2:B:23:TYR:OH	2.09	0.53
1:A:390:ASN:CA	1:A:499:TYR:CE2	2.91	0.53
1:A:393:LEU:HD12	1:A:499:TYR:CD1	2.44	0.53
2:C:25:GLY:HA2	4:C:202:HOH:O	2.10	0.52
1:A:289:VAL:HG22	1:A:296:PHE:HE1	1.75	0.52
1:A:285:SER:HB2	1:A:301:HIS:O	2.09	0.52
1:A:343:ASP:O	1:A:346:ILE:HG13	2.10	0.52
1:A:158:ALA:O	1:A:159:ALA:C	2.47	0.52
1:A:278:GLN:OE1	1:A:334:ARG:HD2	2.10	0.52
1:A:276:PHE:CZ	1:A:334:ARG:HD3	2.45	0.52
1:A:452:TYR:CG	1:A:472:TRP:HH2	2.28	0.52
1:A:160:GLU:C	1:A:162:LEU:H	2.13	0.52
1:A:305:SER:C	1:A:306:ILE:HD12	2.30	0.52
1:A:419:LEU:CD2	4:A:943:HOH:O	2.58	0.52
2:B:87:VAL:HG12	2:B:88:GLN:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:LEU:HD23	1:A:454:LEU:O	2.10	0.51
2:B:57:LEU:HD22	2:B:69:GLN:HE21	1.75	0.51
1:A:506:GLY:O	1:A:507:ASP:CB	2.59	0.51
1:A:275:TYR:HE2	1:A:277:GLU:HG3	1.75	0.51
1:A:302:ARG:HG3	1:A:302:ARG:HH11	1.74	0.51
1:A:419:LEU:O	1:A:421:VAL:CG2	2.58	0.51
2:C:100:THR:N	2:C:104:GLY:O	2.37	0.51
3:A:801:NAG:H82	4:A:946:HOH:O	2.10	0.51
1:A:355:ARG:NH1	1:A:355:ARG:HG3	2.19	0.51
2:C:59:VAL:N	4:C:209:HOH:O	2.04	0.51
1:A:248:GLN:C	1:A:249:LEU:HD13	2.32	0.50
1:A:383:GLN:HA	1:A:386:CYS:HB2	1.92	0.50
1:A:581:HIS:HE1	1:A:596:HIS:HD2	1.59	0.50
2:B:21:CYS:HB2	2:B:36:TRP:HZ2	1.76	0.50
1:A:294:LEU:N	1:A:294:LEU:HD23	2.26	0.50
1:A:334:ARG:NH1	2:C:7:ASP:OD1	2.44	0.50
2:C:25:GLY:CA	4:C:202:HOH:O	2.59	0.50
2:C:33:GLN:OE1	2:C:50:LEU:HD11	2.11	0.50
1:A:391:GLN:HA	1:A:394:CYS:HB3	1.93	0.50
4:A:954:HOH:O	2:B:53:SER:O	2.18	0.50
1:A:156:ASP:CG	1:A:157:VAL:N	2.64	0.50
1:A:487:ILE:O	1:A:488:LYS:O	2.30	0.50
1:A:404:ASP:OD1	1:A:406:ARG:HG3	2.12	0.49
1:A:500:LEU:C	4:A:917:HOH:O	2.49	0.49
1:A:458:PRO:HG2	1:A:526:LEU:HD11	1.94	0.49
1:A:560:LYS:CB	4:A:929:HOH:O	2.56	0.49
1:A:567:GLN:OE1	1:A:582:PHE:HD1	1.95	0.49
2:B:14:GLY:H	2:B:86:ALA:HB3	1.77	0.49
1:A:281:SER:O	1:A:282:ASN:O	2.30	0.49
1:A:549:PHE:HD2	1:A:550:SER:N	2.10	0.49
1:A:390:ASN:OD1	1:A:393:LEU:HB2	2.13	0.49
1:A:430:GLY:HA3	1:A:476:PHE:CE2	2.47	0.49
1:A:508:VAL:HB	1:A:530:ASP:OD1	2.13	0.49
1:A:236:LYS:HD2	1:A:252:HIS:ND1	2.28	0.49
1:A:377:ARG:O	1:A:378:MET:SD	2.71	0.49
2:C:38:ARG:HH12	2:C:90:ASP:HA	1.76	0.49
1:A:531:THR:O	1:A:531:THR:OG1	2.30	0.49
1:A:308:ILE:HG12	1:A:349:LEU:HD12	1.95	0.49
1:A:458:PRO:HB3	2:B:101:PHE:CE2	2.48	0.49
1:A:284:PHE:CE1	4:A:948:HOH:O	2.55	0.48
1:A:352:SER:HB2	1:A:440:GLY:HA2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:488:LYS:C	4:A:958:HOH:O	2.38	0.48
2:C:75:ASN:OD1	2:C:76:PRO:HD2	2.14	0.48
1:A:461:ASN:C	1:A:461:ASN:ND2	2.66	0.48
1:A:320:GLN:OE1	1:A:334:ARG:HD2	2.13	0.48
1:A:464:LEU:CG	2:B:100:THR:O	2.58	0.48
2:C:11:VAL:HG21	2:C:17:ALA:HB2	1.95	0.48
1:A:329:SER:HB2	1:A:330:PRO:HD2	1.96	0.48
1:A:418:SER:O	1:A:419:LEU:CB	2.51	0.48
1:A:253:ARG:HH11	1:A:253:ARG:CG	2.27	0.48
1:A:470:LEU:HD12	1:A:471:GLU:H	1.78	0.47
1:A:421:VAL:HG12	1:A:422:GLU:N	2.30	0.47
1:A:437:HIS:CD2	1:A:439:SER:H	2.23	0.47
1:A:337:VAL:HG13	1:A:423:LEU:O	2.13	0.47
1:A:454:LEU:C	1:A:454:LEU:HD23	2.35	0.47
1:A:488:LYS:HB3	1:A:489:GLU:H	1.56	0.47
1:A:526:LEU:CD2	2:B:102:PRO:HB2	2.45	0.47
2:C:1:GLY:HA3	2:C:24:ARG:O	2.14	0.47
1:A:390:ASN:O	1:A:394:CYS:N	2.40	0.47
1:A:445:LYS:CG	1:A:452:TYR:CE1	2.98	0.47
1:A:472:TRP:CD2	1:A:472:TRP:N	2.83	0.47
2:B:17:ALA:HB3	2:B:83:LEU:HB3	1.97	0.47
1:A:587:ASP:CA	4:A:929:HOH:O	2.62	0.47
2:B:106:PHE:C	2:B:107:GLN:HG3	2.35	0.47
1:A:473:ILE:HA	1:A:474:PRO:HA	1.54	0.47
2:B:87:VAL:HG12	2:B:88:GLN:H	1.79	0.47
1:A:420:THR:O	1:A:420:THR:OG1	2.30	0.47
1:A:340:SER:HB3	1:A:426:LYS:HA	1.97	0.47
1:A:461:ASN:C	1:A:461:ASN:HD22	2.17	0.47
1:A:588:SER:OG	1:A:588:SER:O	2.30	0.47
1:A:445:LYS:HA	1:A:452:TYR:CD1	2.50	0.46
1:A:461:ASN:O	1:A:461:ASN:CG	2.52	0.46
1:A:474:PRO:O	1:A:475:ARG:HG3	2.15	0.46
2:C:50:LEU:HD12	2:C:51:LEU:N	2.30	0.46
1:A:255:PHE:CE1	1:A:322:VAL:HG21	2.50	0.46
2:B:39:VAL:HG21	2:B:93:GLU:HB3	1.97	0.46
2:C:13:LEU:HD21	2:C:88:GLN:NE2	2.30	0.46
1:A:374:ASP:O	1:A:375:LYS:C	2.52	0.46
1:A:525:VAL:CG2	1:A:526:LEU:N	2.78	0.46
1:A:308:ILE:HA	1:A:309:PRO:HD3	1.74	0.46
2:B:66:ARG:CZ	2:B:83:LEU:HD11	2.46	0.46
1:A:285:SER:HA	1:A:302:ARG:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:529:TYR:CG	1:A:563:PRO:HG3	2.49	0.46
2:C:12:VAL:O	2:C:15:GLN:HB3	2.16	0.46
1:A:406:ARG:HH11	1:A:406:ARG:HG3	1.81	0.45
1:A:224:SER:C	1:A:225:GLN:HG2	2.35	0.45
1:A:390:ASN:N	1:A:499:TYR:HD2	2.03	0.45
1:A:355:ARG:NH1	1:A:355:ARG:CG	2.79	0.45
1:A:248:GLN:C	1:A:249:LEU:CD1	2.85	0.45
1:A:276:PHE:CZ	1:A:334:ARG:NH1	2.85	0.45
1:A:362:GLN:OE1	1:A:414:SER:HB2	2.17	0.45
1:A:381:CYS:HA	1:A:490:ALA:CB	2.45	0.45
1:A:232:TYR:CD1	1:A:232:TYR:N	2.84	0.45
3:A:801:NAG:O7	3:A:801:NAG:O3	2.30	0.45
2:B:101:PHE:CG	2:B:102:PRO:HA	2.52	0.45
1:A:337:VAL:HA	1:A:338:PRO:HD3	1.86	0.45
1:A:352:SER:HB2	1:A:353:SER:H	1.64	0.45
2:C:34:VAL:HG21	2:C:79:GLY:HA3	1.99	0.45
1:A:165:ALA:C	1:A:166:LEU:HG	2.38	0.45
1:A:401:PRO:HA	1:A:404:ASP:HB3	1.99	0.44
2:C:38:ARG:CG	2:C:48:LEU:HD11	2.47	0.44
2:C:33:GLN:NE2	2:C:50:LEU:HD11	2.32	0.44
2:C:59:VAL:HG12	2:C:60:SER:N	2.32	0.44
1:A:162:LEU:O	1:A:163:MET:C	2.54	0.44
1:A:216:VAL:CG2	1:A:219:ILE:HD11	2.46	0.44
1:A:212:ARG:HA	1:A:212:ARG:HD3	1.78	0.44
1:A:225:GLN:HB2	1:A:226:GLY:H	1.63	0.44
1:A:418:SER:CB	4:A:938:HOH:O	2.63	0.44
2:B:97:ARG:HG2	2:B:107:GLN:HG2	1.99	0.44
2:B:67:VAL:CG2	2:B:83:LEU:HD13	2.48	0.44
1:A:357:VAL:HG11	1:A:445:LYS:HD2	1.99	0.44
1:A:360:ASP:HB3	3:A:801:NAG:O6	2.18	0.44
1:A:362:GLN:HE21	3:A:801:NAG:H62	1.83	0.44
1:A:192:THR:O	1:A:193:THR:O	2.36	0.44
1:A:445:LYS:HG3	1:A:452:TYR:CE1	2.53	0.44
3:A:801:NAG:C4	4:A:902:HOH:O	2.40	0.44
1:A:455:THR:C	1:A:456:ILE:HG13	2.39	0.43
2:C:19:LEU:HA	2:C:20:PRO:HD3	1.82	0.43
2:B:71:PRO:HA	2:B:72:PRO:HD3	1.91	0.43
1:A:159:ALA:C	1:A:162:LEU:HB2	2.34	0.43
1:A:543:TYR:CD2	1:A:548:SER:HB2	2.53	0.43
1:A:222:MET:SD	1:A:289:VAL:HG13	2.59	0.43
1:A:306:ILE:CG2	1:A:307:THR:N	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:LYS:HE3	1:A:277:GLU:OE1	2.19	0.43
2:C:48:LEU:HD21	2:C:94:TYR:CE1	2.53	0.43
1:A:445:LYS:HG2	1:A:452:TYR:CE1	2.54	0.43
2:C:100:THR:OG1	2:C:104:GLY:O	2.35	0.43
1:A:254:VAL:HG13	1:A:277:GLU:HG2	2.01	0.43
1:A:262:ASN:C	1:A:264:GLY:H	2.23	0.43
1:A:540:TYR:CZ	1:A:579:CYS:HB3	2.54	0.43
1:A:160:GLU:O	1:A:161:GLU:C	2.58	0.43
1:A:157:VAL:HG12	1:A:158:ALA:H	1.82	0.42
1:A:553:TYR:HE1	1:A:600:VAL:HG22	1.84	0.42
2:B:19:LEU:O	2:B:80:SER:HB3	2.19	0.42
1:A:376:LEU:N	1:A:376:LEU:CD2	2.30	0.42
1:A:459:MET:HB3	1:A:463:ALA:HB3	2.00	0.42
1:A:488:LYS:O	1:A:489:GLU:C	2.57	0.42
1:A:500:LEU:CB	2:B:53:SER:OG	2.67	0.42
1:A:553:TYR:HE1	1:A:600:VAL:CG2	2.32	0.42
1:A:437:HIS:HD2	1:A:439:SER:N	2.09	0.42
2:B:15:GLN:HB3	2:B:16:ASP:H	1.56	0.42
1:A:233:LEU:HD21	1:A:253:ARG:NH1	2.35	0.42
1:A:390:ASN:N	1:A:499:TYR:CE2	2.86	0.42
1:A:457:PRO:HB3	1:A:512:SER:O	2.19	0.42
1:A:361:ASN:HD22	1:A:418:SER:HB3	1.81	0.42
1:A:487:ILE:CG2	1:A:487:ILE:O	2.67	0.42
1:A:371:ARG:HH11	1:A:429:SER:CB	2.33	0.42
1:A:536:HIS:HB2	1:A:559:ILE:HB	2.01	0.42
2:B:112:LEU:C	4:B:202:HOH:O	2.52	0.42
1:A:262:ASN:N	1:A:263:PRO:HD3	2.35	0.42
1:A:460:LYS:O	1:A:462:LEU:N	2.53	0.42
2:B:39:VAL:CG2	2:B:93:GLU:HB3	2.50	0.42
1:A:456:ILE:HB	1:A:466:VAL:HG13	2.01	0.42
1:A:229:GLY:HA2	1:A:256:GLU:O	2.19	0.42
1:A:278:GLN:HA	1:A:279:PRO:HD3	1.82	0.42
1:A:388:GLY:HA2	2:B:55:TYR:CE2	2.55	0.42
2:B:3:LEU:HD11	2:B:98:VAL:CG2	2.48	0.42
2:C:35:ALA:O	2:C:96:CYS:HA	2.20	0.42
1:A:253:ARG:NH1	1:A:253:ARG:CG	2.83	0.41
1:A:392:ALA:O	1:A:396:ASN:HB2	2.20	0.41
1:A:275:TYR:CE2	1:A:277:GLU:HG3	2.54	0.41
1:A:361:ASN:HD21	1:A:418:SER:CA	2.33	0.41
1:A:371:ARG:HG3	1:A:409:SER:HB3	2.03	0.41
1:A:472:TRP:CE3	1:A:472:TRP:N	2.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:5:THR:HG23	2:B:6:SER:N	2.34	0.41
2:C:101:PHE:HA	2:C:102:PRO:HA	1.83	0.41
1:A:282:ASN:OD1	4:A:912:HOH:O	2.22	0.41
2:B:19:LEU:HD12	2:B:81:VAL:HG23	2.01	0.41
1:A:343:ASP:HA	1:A:344:PRO:HD2	1.92	0.41
1:A:393:LEU:CD1	1:A:499:TYR:CD1	3.04	0.41
1:A:506:GLY:O	1:A:507:ASP:HB3	2.20	0.41
1:A:542:VAL:HG21	1:A:549:PHE:CZ	2.55	0.41
2:B:41:ALA:O	2:B:43:GLU:HG2	2.20	0.41
2:C:59:VAL:CA	4:C:209:HOH:O	2.63	0.41
2:C:88:GLN:C	2:C:90:ASP:H	2.23	0.41
1:A:299:LEU:HD11	1:A:425:ILE:CD1	2.48	0.41
1:A:382:PHE:CD2	1:A:383:GLN:HG3	2.56	0.41
2:B:101:PHE:CD2	2:B:102:PRO:HA	2.55	0.41
1:A:157:VAL:C	1:A:159:ALA:N	2.74	0.41
1:A:276:PHE:CE1	1:A:334:ARG:NH1	2.89	0.41
2:B:5:THR:OG1	4:B:201:HOH:O	1.86	0.41
1:A:162:LEU:HD22	1:A:162:LEU:HA	1.90	0.41
1:A:485:VAL:HA	1:A:486:PRO:HD3	1.90	0.41
1:A:606:CYS:HB3	4:A:923:HOH:O	2.21	0.41
1:A:430:GLY:HA3	1:A:476:PHE:CD2	2.56	0.41
1:A:358:ILE:HG12	1:A:363:ALA:HB2	2.03	0.40
2:C:68:GLU:OE2	2:C:84:ARG:HD2	2.21	0.40
1:A:299:LEU:HD21	1:A:425:ILE:HD13	2.02	0.40
1:A:354:HIS:ND1	1:A:367:VAL:HG12	2.35	0.40
1:A:370:THR:HB	1:A:406:ARG:NH2	2.36	0.40
1:A:400:ALA:HB3	1:A:401:PRO:HD3	2.03	0.40
2:C:43:GLU:HA	2:C:43:GLU:OE2	2.21	0.40
1:A:165:ALA:O	1:A:166:LEU:HG	2.21	0.40
1:A:166:LEU:CD1	1:A:166:LEU:O	2.67	0.40
1:A:556:ARG:HD2	1:A:556:ARG:HA	1.74	0.40
2:B:59:VAL:CG1	2:B:60:SER:N	2.84	0.40

All (2) 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:534:VAL:CG2	2:B:24:ARG:CZ[3_455]	1.88	0.32
1:A:534:VAL:CG2	2:B:24:ARG:NE[3_455]	1.97	0.23

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	401/473 (85%)	318 (79%)	56 (14%)	27 (7%)	1 7
2	B	112/123 (91%)	94 (84%)	12 (11%)	6 (5%)	2 12
2	C	110/123 (89%)	100 (91%)	8 (7%)	2 (2%)	8 34
All	All	623/719 (87%)	512 (82%)	76 (12%)	35 (6%)	2 11

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	158	ALA
1	A	160	GLU
1	A	161	GLU
1	A	282	ASN
1	A	375	LYS
1	A	461	ASN
1	A	488	LYS
1	A	489	GLU
1	A	535	GLU
1	A	536	HIS
1	A	589	GLU
1	A	193	THR
1	A	376	LEU
1	A	421	VAL
2	B	15	GLN
2	B	104	GLY
2	C	26	ASP
1	A	339	LEU
1	A	557	LEU
1	A	588	SER
1	A	347	ASP
1	A	546	SER
1	A	574	ASP
2	B	21	CYS

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Mol	Chain	Res	Type
2	B	40	ASP
2	B	61	PRO
2	B	62	ALA
1	A	163	MET
1	A	352	SER
1	A	390	ASN
2	C	86	ALA
1	A	159	ALA
1	A	462	LEU
1	A	558	PRO
1	A	491	GLY

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	359/412 (87%)	290 (81%)	69 (19%)	1 6
2	B	91/100 (91%)	78 (86%)	13 (14%)	3 14
2	C	89/100 (89%)	78 (88%)	11 (12%)	4 19
All	All	539/612 (88%)	446 (83%)	93 (17%)	2 9

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	GLU
1	A	161	GLU
1	A	162	LEU
1	A	163	MET
1	A	192	THR
1	A	194	ILE
1	A	195	ARG
1	A	225	GLN
1	A	232	TYR
1	A	248	GLN
1	A	249	LEU

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Mol	Chain	Res	Type
1	A	250	SER
1	A	253	ARG
1	A	257	VAL
1	A	268	PRO
1	A	281	SER
1	A	286	ASN
1	A	287	CYS
1	A	289	VAL
1	A	294	LEU
1	A	302	ARG
1	A	305	SER
1	A	308	ILE
1	A	324	LEU
1	A	346	ILE
1	A	349	LEU
1	A	371	ARG
1	A	374	ASP
1	A	375	LYS
1	A	376	LEU
1	A	377	ARG
1	A	389	LYS
1	A	395	GLU
1	A	413	LEU
1	A	416	ASN
1	A	420	THR
1	A	422	GLU
1	A	424	LYS
1	A	434	LEU
1	A	461	ASN
1	A	466	VAL
1	A	472	TRP
1	A	479	SER
1	A	487	ILE
1	A	489	GLU
1	A	494	CYS
1	A	498	THR
1	A	499	TYR
1	A	500	LEU
1	A	505	ASP
1	A	508	VAL
1	A	514	LEU
1	A	521	ASP

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Mol	Chain	Res	Type
1	A	522	LEU
1	A	526	LEU
1	A	531	THR
1	A	532	SER
1	A	533	ARG
1	A	535	GLU
1	A	546	SER
1	A	547	ARG
1	A	548	SER
1	A	549	PHE
1	A	552	PHE
1	A	576	LYS
1	A	580	ARG
1	A	585	LEU
1	A	589	GLU
1	A	600	VAL
2	B	5	THR
2	B	10	THR
2	B	15	GLN
2	B	24	ARG
2	B	29	GLU
2	B	51	LEU
2	B	52	HIS
2	B	66	ARG
2	B	74	ARG
2	B	91	GLU
2	B	96	CYS
2	B	110	LEU
2	B	114	VAL
2	C	2	GLU
2	C	7	ASP
2	C	13	LEU
2	C	21	CYS
2	C	24	ARG
2	C	47	GLU
2	C	50	LEU
2	C	51	LEU
2	C	74	ARG
2	C	95	GLU
2	C	96	CYS

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

Mol	Chain	Res	Type
1	A	225	GLN
1	A	252	HIS
1	A	301	HIS
1	A	361	ASN
1	A	391	GLN
1	A	437	HIS
1	A	536	HIS
1	A	567	GLN
1	A	596	HIS
2	B	46	GLN
2	B	52	HIS
2	B	69	GLN
2	C	15	GLN
2	C	88	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

There are no monosaccharides in this entry.

5.6 Ligand geometry

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	NAG	A	801	1	14,14,15	0.56	0	17,19,21	0.66	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	801	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	801	NAG	C1-C2-N2-C7
3	A	801	NAG	C4-C5-C6-O6
3	A	801	NAG	O5-C5-C6-O6
3	A	801	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	NAG	15	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	411/473 (86%)	0.04	9 (2%) 62 41	24, 72, 165, 292	0
2	B	114/123 (92%)	0.42	10 (8%) 10 4	51, 119, 200, 234	0
2	C	112/123 (91%)	0.05	4 (3%) 42 22	63, 99, 162, 180	0
All	All	637/719 (88%)	0.11	23 (3%) 42 22	24, 91, 174, 292	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	555	PHE	3.9
2	C	101	PHE	3.8
2	B	85	ASN	3.5
2	B	67	VAL	3.5
2	B	11	VAL	3.2
2	B	16	ASP	3.1
1	A	560	LYS	3.0
2	B	12	VAL	2.9
1	A	606	CYS	2.9
1	A	341	THR	2.7
2	B	17	ALA	2.6
1	A	558	PRO	2.5
2	C	41	ALA	2.4
2	B	48	LEU	2.4
2	B	50	LEU	2.4
2	B	53	SER	2.3
2	C	43	GLU	2.2
2	C	27	SER	2.2
2	B	84	ARG	2.2
1	A	554	PRO	2.2
1	A	530	ASP	2.1
1	A	389	LYS	2.1
1	A	309	PRO	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 monosaccharides 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, 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	B-factors(Å^2)	Q<0.9
3	NAG	A	801	14/15	0.82	0.23	96,111,128,130	0

6.5 Other polymers (i)

There are no such residues in this entry.