



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:40 PM GMT

PDB ID : 3OC0
Title : Structure of human DPP-IV with HTS hit (2S,3S,11bS)-3-butyl-9,10-dimethoxy-1,3,4,6,7,11b-hexahydro-2H-pyrido[2,1-a]isoquinolin-2-ylamine
Authors : Hennig, M.; Stihle, M.; Thoma, R.
Deposited on : 2010-08-09
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

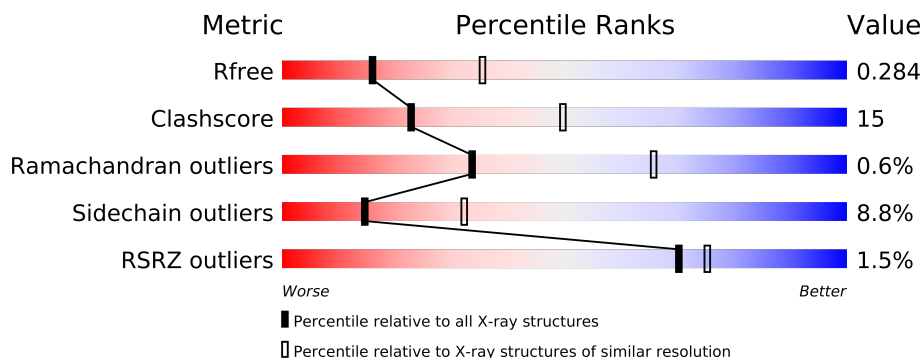
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	728	
1	B	728	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	B	793	-	X
2	NAG	B	797	-	X

2 Entry composition i

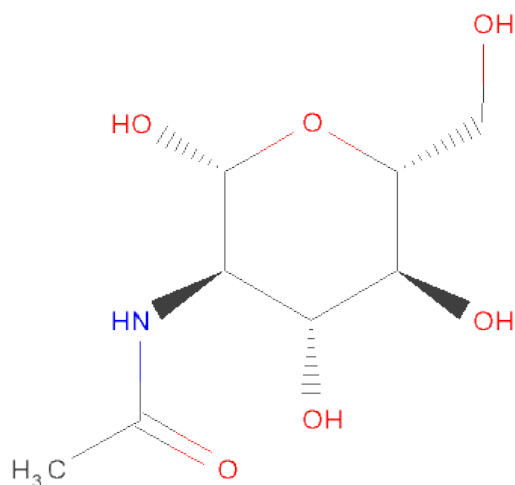
There are 4 unique types of molecules in this entry. The entry contains 12163 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			
1	B	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



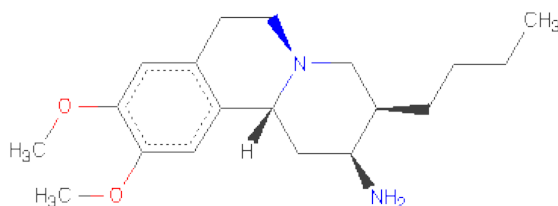
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is (2S,3R,11BR)-3-BUTYL-9,10-DIMETHOXY-1,3,4,6,7,11B-HEXAHYDRO-2H-PYRIDO[2,1-A]ISOQUINOLIN-2-AMINE (three-letter code: B2Q) (formula: $C_{19}H_{30}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			23	19	2	2		
3	B	1	Total	C	N	O	0	0
			23	19	2	2		

- Molecule 4 is water.

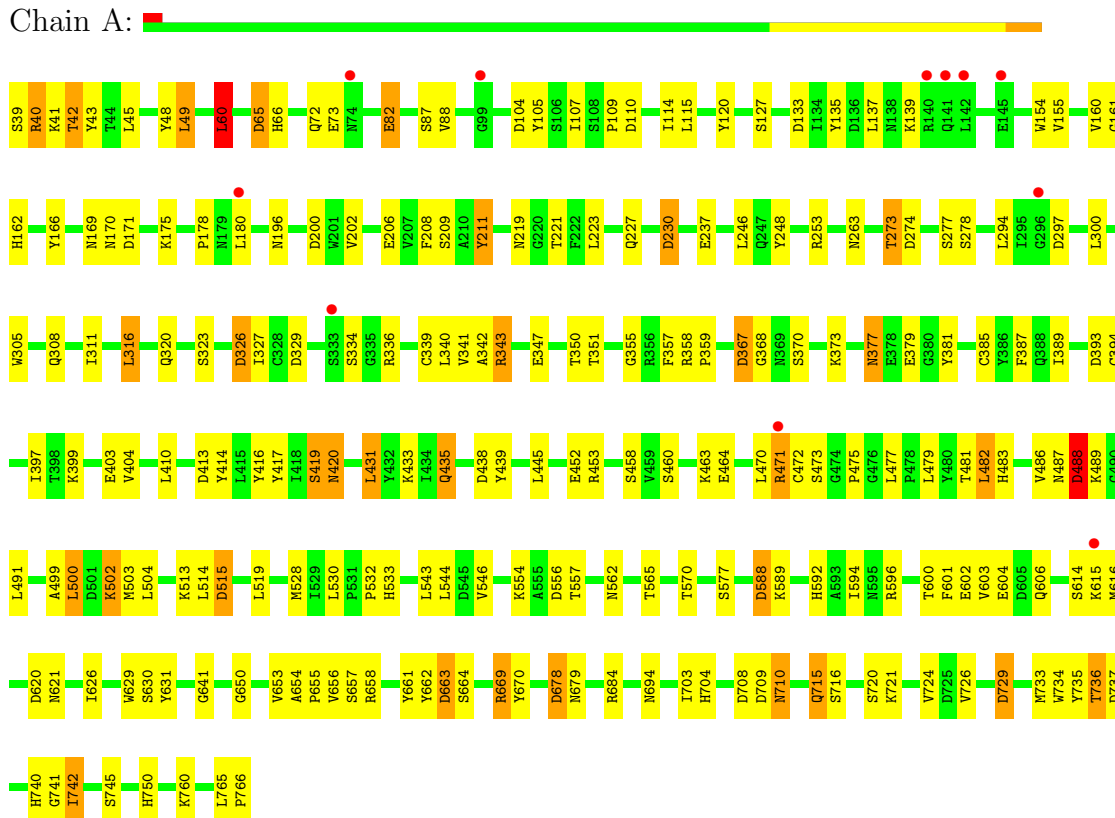
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total	O	0	0
			41	41		
4	B	38	Total	O	0	0
			38	38		

3 Residue-property plots

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.

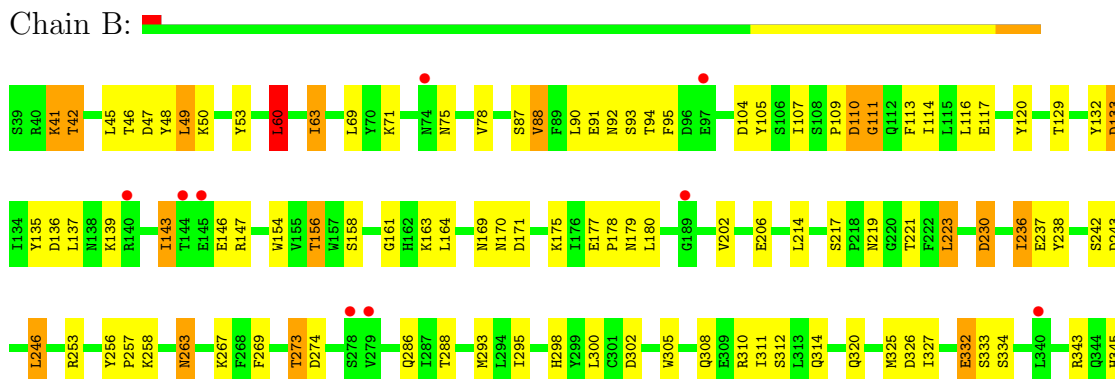
• Molecule 1: Dipeptidyl peptidase 4

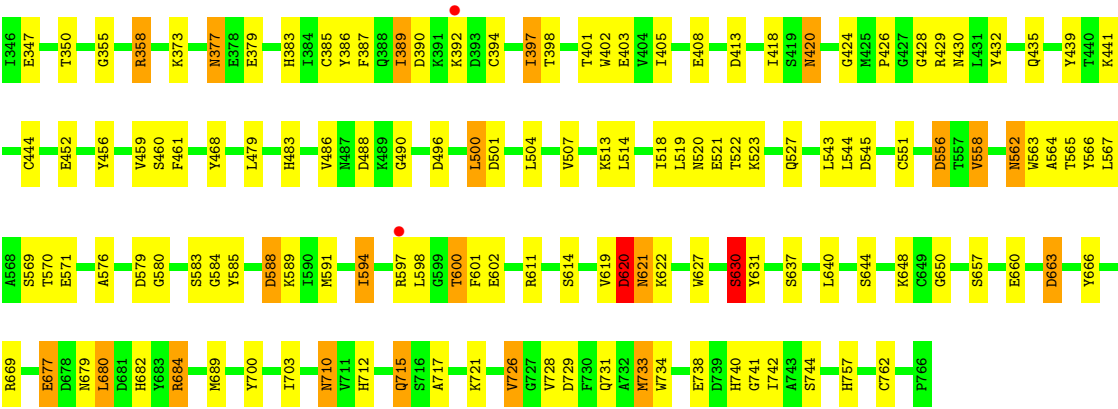
Chain A:



• Molecule 1: Dipeptidyl peptidase 4

Chain B:





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.56Å 68.49Å 421.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.70 10.01 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (12.00-2.70) 71.2 (10.01-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.85 (at 2.71Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.236 , 0.286 0.247 , 0.284	Depositor DCC
R_{free} test set	2042 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	1.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 5.8	EDS
Estimated twinning fraction	0.033 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 40043 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12163	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, B2Q

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	0.76	0/6135	0.94	21/8344 (0.3%)
1	B	0.77	0/6135	0.96	24/8344 (0.3%)
All	All	0.76	0/12270	0.95	45/16688 (0.3%)

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	A	1	0

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	358	ARG	NE-CZ-NH2	-7.51	116.55	120.30
1	A	678	ASP	CB-CG-OD2	7.08	124.67	118.30
1	B	110	ASP	CB-CG-OD2	6.92	124.53	118.30
1	B	729	ASP	CB-CG-OD2	6.88	124.49	118.30
1	B	501	ASP	CB-CG-OD2	6.85	124.47	118.30
1	B	133	ASP	CB-CG-OD2	6.85	124.46	118.30
1	B	663	ASP	CB-CG-OD2	6.74	124.37	118.30
1	A	588	ASP	CB-CG-OD2	6.64	124.28	118.30
1	A	230	ASP	CB-CG-OD2	6.62	124.26	118.30
1	A	620	ASP	CB-CG-OD2	6.61	124.25	118.30
1	B	274	ASP	CB-CG-OD2	6.61	124.25	118.30
1	B	579	ASP	CB-CG-OD2	6.46	124.12	118.30
1	A	297	ASP	CB-CG-OD2	6.35	124.02	118.30
1	A	737	ASP	CB-CG-OD2	6.29	123.96	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	367	ASP	CB-CG-OD2	6.28	123.95	118.30
1	B	588	ASP	CB-CG-OD2	6.23	123.90	118.30
1	B	390	ASP	CB-CG-OD2	6.19	123.87	118.30
1	B	111	GLY	N-CA-C	6.13	128.43	113.10
1	A	171	ASP	CB-CG-OD2	6.10	123.79	118.30
1	A	329	ASP	CB-CG-OD2	5.88	123.59	118.30
1	B	620	ASP	CB-CG-OD2	5.86	123.58	118.30
1	A	515	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	133	ASP	CB-CG-OD2	5.77	123.49	118.30
1	B	243	ASP	CB-CG-OD2	5.65	123.39	118.30
1	B	47	ASP	CB-CG-OD2	5.45	123.20	118.30
1	A	413	ASP	CB-CG-OD2	5.44	123.20	118.30
1	B	230	ASP	CB-CG-OD2	5.42	123.18	118.30
1	B	721	LYS	CD-CE-NZ	5.41	124.13	111.70
1	A	488	ASP	CB-CG-OD2	5.38	123.14	118.30
1	A	200	ASP	CB-CG-OD2	5.35	123.12	118.30
1	A	60	LEU	CA-CB-CG	5.35	127.60	115.30
1	B	60	LEU	CA-CB-CG	5.34	127.58	115.30
1	A	104	ASP	CB-CG-OD2	5.33	123.10	118.30
1	B	171	ASP	CB-CG-OD2	5.33	123.09	118.30
1	B	104	ASP	CB-CG-OD2	5.33	123.09	118.30
1	B	496	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	488	ASP	CB-CG-OD2	5.28	123.06	118.30
1	B	358	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	A	709	ASP	CB-CG-OD2	5.25	123.02	118.30
1	B	413	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	65	ASP	CB-CG-OD2	5.16	122.95	118.30
1	B	556	ASP	CB-CG-OD2	5.12	122.91	118.30
1	A	274	ASP	CB-CG-OD2	5.04	122.84	118.30
1	A	729	ASP	CB-CG-OD2	5.03	122.82	118.30
1	A	40	ARG	N-CA-C	5.01	124.52	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	40	ARG	CA

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5963	0	5681	146	0
1	B	5963	0	5680	198	0
2	A	56	0	52	2	0
2	B	56	0	52	3	0
3	A	23	0	30	2	0
3	B	23	0	30	4	0
4	A	41	0	0	5	0
4	B	38	0	0	13	0
All	All	12163	0	11525	346	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (346) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:600:THR:HG22	1:B:601:PHE:H	1.13	1.12
1:B:267:LYS:HE3	4:B:772:HOH:O	1.56	1.06
1:B:543:LEU:HD21	1:B:627:TRP:HD1	1.22	1.02
1:A:347:GLU:OE2	1:A:373:LYS:NZ	2.01	0.94
1:B:600:THR:HG22	1:B:601:PHE:N	1.83	0.93
1:B:221:THR:O	1:B:273:THR:HB	1.71	0.91
1:A:377:ASN:C	1:A:377:ASN:HD22	1.74	0.90
1:A:160:VAL:HG12	1:A:160:VAL:O	1.68	0.90
1:B:377:ASN:HD22	1:B:377:ASN:C	1.74	0.89
1:B:620:ASP:OD1	1:B:620:ASP:C	2.10	0.88
1:B:267:LYS:CE	4:B:772:HOH:O	2.14	0.87
1:B:680:LEU:HD22	1:B:684:ARG:HD3	1.58	0.86
1:A:704:HIS:HD1	1:A:716:SER:HG	0.87	0.84
1:A:735:TYR:OH	1:A:750:HIS:HD2	1.60	0.84
1:B:253:ARG:HD2	4:B:767:HOH:O	1.77	0.84
1:B:87:SER:OG	2:B:794:NAG:O7	1.97	0.81
1:B:267:LYS:HD2	1:B:286:GLN:NE2	1.95	0.80
1:B:117:GLU:HG3	1:B:132:TYR:CE1	2.17	0.79
1:B:75:ASN:OD1	1:B:92:ASN:HB3	1.81	0.79

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:657:SER:H	1:B:715:GLN:NE2	1.81	0.79
1:B:320:GLN:OE1	1:B:669:ARG:HD3	1.83	0.78
3:B:900:B2Q:HH	4:B:11:HOH:O	1.81	0.78
1:B:347:GLU:OE2	1:B:373:LYS:NZ	2.14	0.78
1:A:596:ARG:NH2	1:A:678:ASP:OD1	2.16	0.78
1:B:46:THR:HG22	1:B:50:LYS:HD3	1.66	0.78
1:B:731:GLN:NE2	4:B:779:HOH:O	2.16	0.78
1:B:600:THR:CG2	1:B:601:PHE:H	1.96	0.77
1:B:562:ASN:HD22	1:B:562:ASN:C	1.90	0.75
1:A:487:ASN:O	1:A:489:LYS:N	2.20	0.74
1:A:403:GLU:H	1:A:420:ASN:HD21	1.33	0.74
1:B:175:LYS:NZ	1:B:178:PRO:O	2.21	0.74
1:B:600:THR:CG2	1:B:601:PHE:N	2.49	0.73
1:A:305:TRP:CZ3	1:A:311:ILE:HG12	2.23	0.73
1:B:543:LEU:HD21	1:B:627:TRP:CD1	2.15	0.72
1:B:621:ASN:HD22	1:B:622:LYS:N	1.88	0.72
1:A:414:TYR:CE1	1:A:435:GLN:HG3	2.24	0.72
1:B:242:SER:OG	1:B:246:LEU:HD12	1.88	0.72
1:A:109:PRO:HD2	1:A:161:GLY:O	1.91	0.71
1:B:236:ILE:HD12	1:B:712:HIS:CD2	2.26	0.70
1:B:163:LYS:HZ3	1:B:273:THR:HG22	1.54	0.70
1:B:401:THR:HG22	1:B:401:THR:O	1.90	0.70
1:B:432:TYR:CE2	1:B:444:CYS:HB2	2.27	0.69
2:A:793:NAG:H3	2:A:793:NAG:H83	1.75	0.68
1:B:377:ASN:C	1:B:377:ASN:ND2	2.47	0.68
1:B:403:GLU:H	1:B:420:ASN:HD21	1.41	0.68
1:A:357:PHE:CE1	3:A:900:B2Q:HGA	2.29	0.68
1:B:598:LEU:O	1:B:682:HIS:HE1	1.76	0.68
1:A:514:LEU:HD23	1:A:514:LEU:C	2.15	0.67
1:B:403:GLU:OE1	1:B:585:TYR:HA	1.95	0.67
1:A:160:VAL:CG1	1:A:160:VAL:O	2.40	0.67
1:B:345:HIS:CE1	4:B:28:HOH:O	2.46	0.67
1:A:657:SER:H	1:A:715:GLN:NE2	1.93	0.66
1:B:219:ASN:N	1:B:308:GLN:OE1	2.29	0.66
1:B:620:ASP:OD1	1:B:622:LYS:N	2.29	0.66
1:A:221:THR:O	1:A:273:THR:HB	1.95	0.66
1:A:482:LEU:HD13	1:A:491:LEU:HD12	1.79	0.65
1:B:518:ILE:O	1:B:519:LEU:HD12	1.96	0.65
1:A:397:ILE:HG22	1:A:439:TYR:CE2	2.31	0.65
1:A:735:TYR:OH	1:A:750:HIS:CD2	2.48	0.65
1:A:594:ILE:CD1	1:A:601:PHE:HB2	2.28	0.64
1:A:42:THR:CG2	1:A:570:THR:OG1	2.46	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:377:ASN:ND2	1:A:377:ASN:C	2.44	0.64
1:B:680:LEU:CD2	1:B:684:ARG:HD3	2.26	0.63
1:B:237:GLU:HG2	1:B:253:ARG:HG2	1.81	0.63
1:B:63:ILE:HD13	1:B:69:LEU:HG	1.80	0.63
1:B:163:LYS:NZ	1:B:273:THR:HG22	2.14	0.62
1:B:267:LYS:HD2	1:B:286:GLN:HE22	1.62	0.62
1:B:42:THR:HG23	1:B:570:THR:OG1	2.00	0.62
1:B:621:ASN:C	1:B:621:ASN:HD22	2.02	0.61
1:B:267:LYS:HG3	4:B:772:HOH:O	1.99	0.61
1:B:660:GLU:OE2	4:B:778:HOH:O	2.16	0.61
1:B:305:TRP:CZ3	1:B:311:ILE:HG12	2.36	0.61
1:A:273:THR:O	1:A:273:THR:HG23	2.00	0.60
1:A:351:THR:OG1	1:A:592:HIS:HD2	1.83	0.60
1:B:428:GLY:O	1:B:429:ARG:HG2	2.02	0.60
1:A:736:THR:HG21	1:B:717:ALA:O	2.02	0.60
1:B:75:ASN:OD1	1:B:92:ASN:CB	2.49	0.60
1:B:137:LEU:C	1:B:139:LYS:H	2.04	0.60
1:B:543:LEU:HD23	1:B:544:LEU:N	2.17	0.59
1:B:327:ILE:HD13	1:B:389:ILE:HG12	1.84	0.59
1:A:453:ARG:O	1:A:475:PRO:HD2	2.03	0.59
1:B:594:ILE:HD11	1:B:602:GLU:N	2.17	0.59
1:B:435:GLN:NE2	1:B:441:LYS:HD3	2.18	0.58
1:B:269:PHE:CE1	1:B:286:GLN:HB2	2.38	0.58
1:B:611:ARG:O	1:B:614:SER:HB2	2.02	0.58
1:B:179:ASN:OD1	1:B:180:LEU:HD22	2.02	0.58
1:A:206:GLU:OE2	1:A:663:ASP:OD2	2.19	0.58
1:B:136:ASP:O	1:B:139:LYS:O	2.22	0.58
1:B:42:THR:CG2	1:B:570:THR:OG1	2.51	0.58
1:B:63:ILE:CD1	1:B:69:LEU:CD1	2.81	0.57
1:B:163:LYS:NZ	1:B:273:THR:CG2	2.67	0.57
1:A:305:TRP:CE3	1:A:311:ILE:HG12	2.39	0.57
1:B:242:SER:HG	1:B:246:LEU:HD12	1.68	0.57
1:A:340:LEU:O	1:A:342:ALA:N	2.37	0.57
1:B:257:PRO:CB	1:B:263:ASN:HD22	2.18	0.57
1:B:69:LEU:CD1	1:B:107:ILE:HD12	2.35	0.56
1:A:470:LEU:HD12	1:A:483:HIS:CE1	2.41	0.56
1:A:72:GLN:O	1:A:73:GLU:HB2	2.04	0.56
1:A:115:LEU:HD11	1:A:155:VAL:HG11	1.87	0.56
1:B:657:SER:H	1:B:715:GLN:HE21	1.52	0.56
1:B:598:LEU:O	1:B:682:HIS:CE1	2.58	0.56
1:B:257:PRO:HB2	1:B:263:ASN:ND2	2.21	0.56
1:B:461:PHE:CD1	1:B:468:TYR:HB3	2.41	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:416:TYR:CE2	1:A:433:LYS:HG3	2.41	0.56
1:A:602:GLU:OE2	1:A:631:TYR:OH	2.17	0.55
1:A:114:ILE:HG23	1:A:135:TYR:HB3	1.87	0.55
1:A:614:SER:O	1:A:615:LYS:C	2.44	0.55
1:A:470:LEU:N	1:A:481:THR:O	2.38	0.55
1:A:326:ASP:OD2	1:A:339:CYS:HB3	2.06	0.55
1:A:760:LYS:HE2	4:A:38:HOH:O	2.06	0.55
1:B:710:ASN:C	1:B:710:ASN:HD22	2.09	0.55
1:B:397:ILE:HG13	1:B:439:TYR:CE2	2.42	0.54
3:B:900:B2Q:CH	4:B:11:HOH:O	2.49	0.54
1:A:237:GLU:HG2	1:A:253:ARG:HG2	1.90	0.54
1:B:105:TYR:CD1	1:B:105:TYR:C	2.80	0.54
1:B:703:ILE:HA	1:B:733:MET:O	2.07	0.54
1:B:387:PHE:CE2	1:B:394:CYS:HB3	2.43	0.54
1:B:588:ASP:O	1:B:589:LYS:C	2.44	0.54
1:A:721:LYS:NZ	1:B:242:SER:O	2.39	0.54
1:B:332:GLU:HG2	1:B:333:SER:N	2.23	0.53
1:B:520:ASN:O	1:B:521:GLU:HB2	2.08	0.53
1:B:562:ASN:ND2	1:B:562:ASN:C	2.60	0.53
1:B:69:LEU:HD13	1:B:107:ILE:HD12	1.90	0.53
1:B:257:PRO:CB	1:B:263:ASN:ND2	2.72	0.53
1:B:424:GLY:O	1:B:426:PRO:HD3	2.08	0.53
1:B:680:LEU:O	1:B:680:LEU:HD22	2.09	0.53
1:B:620:ASP:O	1:B:620:ASP:OD1	2.25	0.53
1:A:500:LEU:HA	1:A:503:MET:CE	2.39	0.53
1:A:355:GLY:HA3	1:A:358:ARG:O	2.08	0.53
1:B:387:PHE:CD2	1:B:394:CYS:HB3	2.43	0.53
1:A:82:GLU:OE1	1:A:82:GLU:HA	2.07	0.53
1:A:641:GLY:O	1:A:694:ASN:ND2	2.36	0.53
1:A:334:SER:OG	1:A:336:ARG:HG3	2.09	0.53
1:A:403:GLU:H	1:A:420:ASN:ND2	2.05	0.53
1:B:594:ILE:HG23	1:B:594:ILE:O	2.08	0.53
2:A:793:NAG:H3	2:A:793:NAG:C8	2.37	0.53
1:B:405:ILE:HD13	1:B:429:ARG:CD	2.38	0.53
1:A:703:ILE:HA	1:A:733:MET:O	2.09	0.53
1:A:273:THR:O	1:A:273:THR:CG2	2.57	0.52
1:A:41:LYS:HG2	1:A:42:THR:N	2.24	0.52
1:A:514:LEU:HD23	1:A:515:ASP:N	2.24	0.52
1:A:657:SER:H	1:A:715:GLN:HE21	1.56	0.52
1:B:377:ASN:ND2	1:B:379:GLU:H	2.07	0.52
1:B:386:TYR:HB2	1:B:397:ILE:CD1	2.40	0.52
1:B:594:ILE:HD11	1:B:602:GLU:H	1.73	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:267:LYS:NZ	1:B:286:GLN:HE22	2.07	0.52
1:B:435:GLN:NE2	1:B:441:LYS:CD	2.73	0.51
1:A:105:TYR:HA	1:A:115:LEU:O	2.10	0.51
1:B:377:ASN:ND2	1:B:379:GLU:N	2.57	0.51
1:A:514:LEU:CD2	1:A:514:LEU:C	2.78	0.51
1:A:377:ASN:ND2	1:A:379:GLU:H	2.08	0.51
1:A:463:LYS:O	1:A:464:GLU:HB2	2.09	0.51
1:A:654:ALA:N	1:A:655:PRO:HD3	2.27	0.50
1:A:73:GLU:N	4:A:774:HOH:O	2.39	0.50
1:A:43:TYR:CD2	1:A:565:THR:HG22	2.46	0.50
1:A:196:ASN:OD1	1:A:227:GLN:HG3	2.10	0.50
1:A:528:MET:CE	1:A:530:LEU:HD21	2.42	0.50
1:B:397:ILE:HG13	1:B:439:TYR:CD2	2.47	0.50
1:A:114:ILE:CG2	1:A:135:TYR:HB3	2.41	0.50
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.93	0.50
1:A:720:SER:O	1:A:724:VAL:HG23	2.11	0.50
1:B:429:ARG:HG3	1:B:456:TYR:CZ	2.47	0.49
1:B:666:TYR:CD2	3:B:900:B2Q:HR	2.47	0.49
1:A:458:SER:OG	1:A:471:ARG:CG	2.60	0.49
1:B:513:LYS:O	1:B:527:GLN:HA	2.12	0.49
1:A:596:ARG:HA	1:A:670:TYR:O	2.13	0.49
1:A:499:ALA:O	1:A:502:LYS:HG3	2.13	0.49
1:B:45:LEU:HG	1:B:49:LEU:HD22	1.95	0.49
1:B:110:ASP:CG	1:B:161:GLY:H	2.16	0.49
1:B:137:LEU:C	1:B:139:LYS:N	2.66	0.49
1:A:594:ILE:HD11	1:A:602:GLU:H	1.78	0.48
1:B:420:ASN:C	1:B:420:ASN:HD22	2.16	0.48
1:B:41:LYS:HE3	1:B:53:TYR:OH	2.13	0.48
1:B:483:HIS:CD2	1:B:490:GLY:HA2	2.48	0.48
1:B:267:LYS:CG	4:B:772:HOH:O	2.59	0.48
1:B:459:VAL:HG22	1:B:460:SER:N	2.29	0.48
1:A:532:PRO:O	1:A:533:HIS:HB2	2.13	0.48
1:A:60:LEU:C	1:A:60:LEU:HD23	2.34	0.48
1:B:402:TRP:HB2	1:B:420:ASN:ND2	2.28	0.48
1:A:66:HIS:ND1	4:A:781:HOH:O	2.35	0.48
1:A:397:ILE:HG22	1:A:439:TYR:CD2	2.48	0.48
1:A:42:THR:HG21	1:A:570:THR:OG1	2.13	0.48
1:A:734:TRP:HE1	1:A:736:THR:HG22	1.79	0.48
1:B:500:LEU:HD22	1:B:504:LEU:HG	1.96	0.47
1:B:408:GLU:OE1	4:B:24:HOH:O	2.20	0.47
1:B:163:LYS:HZ3	1:B:273:THR:CG2	2.22	0.47
1:B:46:THR:CG2	1:B:50:LYS:HD3	2.41	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:95:PHE:CZ	1:B:116:LEU:HD11	2.50	0.47
2:B:793:NAG:H3	2:B:793:NAG:C8	2.44	0.47
1:A:741:GLY:O	1:A:742:ILE:C	2.53	0.47
1:A:654:ALA:N	1:A:655:PRO:CD	2.77	0.47
1:A:458:SER:OG	1:A:471:ARG:HG3	2.15	0.47
1:B:689:MET:SD	1:B:689:MET:N	2.87	0.47
1:A:629:TRP:HA	1:A:653:VAL:O	2.15	0.47
1:A:419:SER:OG	1:A:420:ASN:N	2.45	0.47
1:A:105:TYR:C	1:A:105:TYR:CD1	2.88	0.47
1:A:615:LYS:O	1:A:616:MET:C	2.53	0.47
1:A:404:VAL:HG13	1:A:417:TYR:CD2	2.49	0.47
1:B:88:VAL:O	1:B:88:VAL:HG22	2.14	0.47
1:B:562:ASN:ND2	1:B:565:THR:H	2.12	0.47
1:A:656:VAL:HG21	3:A:900:B2Q:HP	1.96	0.47
1:B:92:ASN:OD1	1:B:93:SER:N	2.48	0.47
1:B:564:ALA:O	1:B:565:THR:C	2.52	0.47
1:A:137:LEU:O	1:A:139:LYS:O	2.33	0.47
1:A:323:SER:OG	1:A:347:GLU:HB2	2.14	0.46
1:B:405:ILE:HD13	1:B:429:ARG:NE	2.30	0.46
1:A:596:ARG:N	1:A:670:TYR:O	2.48	0.46
1:A:340:LEU:HB3	1:A:343:ARG:HD2	1.97	0.46
1:B:543:LEU:HD23	1:B:543:LEU:C	2.35	0.46
1:B:435:GLN:HE22	1:B:441:LYS:HD3	1.79	0.46
1:B:677:GLU:H	1:B:677:GLU:HG3	1.45	0.46
1:B:164:LEU:HB2	1:B:175:LYS:HB2	1.96	0.46
1:B:648:LYS:HE3	1:B:762:CYS:O	2.15	0.46
1:A:316:LEU:HD22	1:A:320:GLN:HA	1.98	0.46
1:B:343:ARG:O	1:B:343:ARG:HG2	2.16	0.46
1:B:48:TYR:CD2	1:B:49:LEU:HD13	2.51	0.46
1:B:562:ASN:HD21	1:B:564:ALA:HB3	1.80	0.46
1:B:598:LEU:HD22	1:B:631:TYR:OH	2.16	0.46
1:B:741:GLY:O	1:B:742:ILE:C	2.53	0.46
1:A:127:SER:HB3	1:A:211:TYR:CD1	2.51	0.46
1:B:584:GLY:O	1:B:585:TYR:HB2	2.15	0.46
1:B:435:GLN:HE21	1:B:441:LYS:HB3	1.81	0.46
1:A:300:LEU:HD23	1:A:300:LEU:C	2.36	0.46
1:A:305:TRP:CZ3	1:A:311:ILE:CG1	2.97	0.46
1:B:405:ILE:HD13	1:B:429:ARG:HD3	1.97	0.46
1:A:546:VAL:HG22	1:A:606:GLN:OE1	2.16	0.46
1:B:42:THR:OG1	1:B:569:SER:O	2.34	0.45
1:A:588:ASP:O	1:A:589:LYS:C	2.53	0.45
1:A:726:VAL:O	1:A:726:VAL:CG1	2.64	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:345:HIS:HE1	4:B:28:HOH:O	1.91	0.45
1:A:154:TRP:O	1:A:166:TYR:HA	2.16	0.45
1:A:219:ASN:N	1:A:308:GLN:OE1	2.49	0.45
1:B:734:TRP:C	1:B:734:TRP:CD1	2.90	0.45
1:B:206:GLU:O	3:B:900:B2Q:CA	2.65	0.44
1:A:420:ASN:C	1:A:420:ASN:HD22	2.18	0.44
1:B:403:GLU:H	1:B:420:ASN:ND2	2.13	0.44
1:A:514:LEU:CD1	1:A:557:THR:HG22	2.47	0.44
1:A:500:LEU:HD11	1:A:504:LEU:HD11	1.99	0.44
1:A:710:ASN:C	1:A:710:ASN:HD22	2.21	0.44
1:A:528:MET:HE2	1:A:530:LEU:HD21	1.99	0.44
1:B:137:LEU:O	1:B:139:LYS:N	2.50	0.44
1:A:600:THR:O	1:A:603:VAL:N	2.49	0.44
1:B:133:ASP:OD1	1:B:147:ARG:NH1	2.50	0.44
1:A:482:LEU:CD1	1:A:491:LEU:HD12	2.46	0.44
1:A:387:PHE:CE2	1:A:394:CYS:HB3	2.52	0.44
1:B:117:GLU:HG3	1:B:132:TYR:CZ	2.52	0.44
1:A:562:ASN:O	1:A:565:THR:HB	2.18	0.44
1:B:63:ILE:CD1	1:B:69:LEU:HG	2.46	0.44
1:B:556:ASP:OD1	1:B:558:VAL:HG13	2.18	0.44
1:B:293:MET:O	1:B:298:HIS:HD2	2.01	0.44
1:A:487:ASN:O	1:A:488:ASP:C	2.55	0.44
1:B:169:ASN:O	1:B:170:ASN:HB2	2.17	0.44
1:A:760:LYS:CE	4:A:38:HOH:O	2.65	0.44
1:A:626:ILE:O	1:A:650:GLY:HA2	2.18	0.44
1:B:383:HIS:HD2	1:B:398:THR:HB	1.83	0.44
1:A:669:ARG:HG3	1:A:669:ARG:HH11	1.82	0.44
1:B:267:LYS:CD	1:B:286:GLN:HE22	2.29	0.43
1:B:236:ILE:HD12	1:B:712:HIS:NE2	2.33	0.43
1:A:603:VAL:O	1:A:604:GLU:C	2.56	0.43
1:A:556:ASP:OD1	1:A:556:ASP:C	2.57	0.43
1:B:543:LEU:CD2	1:B:543:LEU:C	2.86	0.43
1:B:562:ASN:O	1:B:563:TRP:C	2.55	0.43
1:B:630:SER:OG	1:B:740:HIS:NE2	2.42	0.43
1:B:177:GLU:HB2	1:B:180:LEU:HD23	2.00	0.43
1:A:658:ARG:HB3	1:A:661:TYR:CD2	2.53	0.43
1:B:355:GLY:HA3	1:B:358:ARG:O	2.18	0.43
1:B:49:LEU:HA	1:B:49:LEU:HD12	1.74	0.43
1:B:738:GLU:OE2	1:B:744:SER:OG	2.24	0.43
1:B:405:ILE:CD1	1:B:429:ARG:HD3	2.48	0.43
1:A:477:LEU:HD22	1:A:500:LEU:HD12	1.99	0.43
1:B:114:ILE:CG2	1:B:135:TYR:HB3	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:158:SER:OG	1:B:163:LYS:HB2	2.18	0.43
1:B:253:ARG:CD	4:B:767:HOH:O	2.50	0.43
1:A:594:ILE:HG23	1:A:594:ILE:O	2.19	0.43
1:A:519:LEU:HA	1:A:519:LEU:HD23	1.64	0.43
1:B:742:ILE:HG22	1:B:742:ILE:O	2.18	0.43
1:A:370:SER:HB2	1:A:387:PHE:O	2.18	0.43
1:B:551:CYS:HB2	1:B:591:MET:SD	2.59	0.43
1:A:729:ASP:OD2	1:B:757:HIS:HD2	2.00	0.43
1:B:597:ARG:O	1:B:600:THR:HB	2.19	0.43
1:B:256:TYR:CZ	1:B:663:ASP:HB3	2.54	0.43
1:B:312:SER:OG	1:B:325:MET:HE3	2.19	0.43
1:A:327:ILE:HB	1:A:343:ARG:HG3	2.01	0.42
1:B:63:ILE:HD13	1:B:69:LEU:CG	2.46	0.42
1:B:726:VAL:HG12	1:B:728:VAL:HG23	2.01	0.42
1:A:377:ASN:ND2	1:A:381:TYR:H	2.16	0.42
1:A:327:ILE:HD13	1:A:389:ILE:HG12	2.00	0.42
1:B:428:GLY:C	1:B:429:ARG:HG2	2.39	0.42
1:B:522:THR:HG22	1:B:523:LYS:N	2.33	0.42
1:B:567:LEU:O	1:B:571:GLU:HB2	2.20	0.42
1:B:305:TRP:CE3	1:B:311:ILE:HG12	2.55	0.42
1:B:320:GLN:OE1	1:B:669:ARG:CD	2.62	0.42
1:A:110:ASP:OD1	1:A:110:ASP:C	2.58	0.42
1:B:594:ILE:CD1	1:B:601:PHE:HB2	2.49	0.42
1:B:562:ASN:HD22	1:B:565:THR:H	1.66	0.42
1:A:765:LEU:HA	1:A:766:PRO:HD3	1.87	0.42
1:A:367:ASP:OD1	1:A:368:GLY:N	2.52	0.42
1:B:143:ILE:HG13	1:B:143:ILE:O	2.13	0.41
1:B:146:GLU:O	1:B:175:LYS:NZ	2.50	0.41
1:B:69:LEU:HD11	1:B:107:ILE:HD12	2.02	0.41
1:A:662:TYR:HE1	1:A:710:ASN:HD22	1.68	0.41
1:A:669:ARG:HG3	1:A:669:ARG:NH1	2.35	0.41
1:A:45:LEU:HG	1:A:49:LEU:HD22	2.02	0.41
1:B:594:ILE:HD12	1:B:601:PHE:HB2	2.01	0.41
1:A:358:ARG:HB2	1:A:359:PRO:HD2	2.02	0.41
1:B:543:LEU:CD2	1:B:627:TRP:HD1	2.11	0.41
1:A:42:THR:HG23	1:A:43:TYR:N	2.35	0.41
1:B:69:LEU:HD23	1:B:78:VAL:HG22	2.03	0.41
1:B:310:ARG:NH2	1:B:389:ILE:HD13	2.35	0.41
1:A:340:LEU:C	1:A:342:ALA:N	2.70	0.41
1:A:500:LEU:HD22	1:A:504:LEU:HG	2.02	0.41
1:A:554:LYS:HB3	1:A:577:SER:HB3	2.02	0.41
1:A:65:ASP:N	1:A:65:ASP:OD1	2.52	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:708:ASP:OD2	1:A:740:HIS:HA	2.21	0.41
1:B:630:SER:O	1:B:631:TYR:C	2.59	0.41
2:B:793:NAG:H82	2:B:793:NAG:C1	2.51	0.41
1:B:154:TRP:NE1	1:B:156:THR:OG1	2.53	0.41
1:B:113:PHE:CE1	1:B:178:PRO:HG2	2.56	0.41
1:B:236:ILE:HD13	1:B:238:TYR:HD1	1.86	0.41
1:A:48:TYR:CD2	1:A:49:LEU:HD13	2.55	0.41
1:B:223:LEU:HD23	1:B:223:LEU:HA	1.85	0.41
1:A:431:LEU:HD13	1:A:445:LEU:HD12	2.01	0.41
1:A:248:TYR:CZ	1:B:258:LYS:HD2	2.55	0.41
1:A:340:LEU:C	1:A:342:ALA:H	2.24	0.41
1:B:397:ILE:CG1	1:B:439:TYR:CE2	3.03	0.41
1:B:580:GLY:O	1:B:583:SER:OG	2.26	0.41
1:A:500:LEU:HA	1:A:503:MET:HE2	2.01	0.41
1:A:662:TYR:HE1	1:A:710:ASN:ND2	2.19	0.41
1:A:208:PHE:O	1:A:209:SER:C	2.56	0.41
1:B:60:LEU:C	1:B:60:LEU:CD2	2.89	0.41
1:B:120:TYR:CD1	1:B:120:TYR:C	2.94	0.41
1:A:420:ASN:C	1:A:420:ASN:ND2	2.75	0.41
1:A:263:ASN:HD21	1:A:664:SER:CB	2.34	0.41
1:B:109:PRO:HD2	1:B:161:GLY:O	2.21	0.40
1:B:637:SER:HG	1:B:700:TYR:HH	1.69	0.40
1:B:545:ASP:N	1:B:576:ALA:O	2.54	0.40
1:B:397:ILE:HG12	1:B:439:TYR:CE1	2.56	0.40
1:A:175:LYS:NZ	1:A:178:PRO:O	2.54	0.40
1:B:267:LYS:CE	1:B:286:GLN:HE22	2.34	0.40
1:A:110:ASP:OD2	1:A:162:HIS:ND1	2.53	0.40
1:B:418:ILE:HA	1:B:430:ASN:O	2.20	0.40
1:A:73:GLU:CA	4:A:774:HOH:O	2.68	0.40
1:A:169:ASN:O	1:A:170:ASN:HB2	2.22	0.40
1:A:105:TYR:HB2	1:A:114:ILE:HD11	2.04	0.40
1:B:500:LEU:HD22	1:B:504:LEU:CD1	2.52	0.40
1:B:114:ILE:O	1:B:114:ILE:HG23	2.21	0.40

There are no symmetry-related clashes.

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	726/728 (100%)	676 (93%)	44 (6%)	6 (1%)	27	58
1	B	726/728 (100%)	674 (93%)	50 (7%)	2 (0%)	50	82
All	All	1452/1456 (100%)	1350 (93%)	94 (6%)	8 (1%)	33	66

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	438	ASP
1	A	488	ASP
1	B	111	GLY
1	A	393	ASP
1	B	630	SER
1	A	742	ILE
1	A	341	VAL
1	A	486	VAL

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	653/653 (100%)	599 (92%)	54 (8%)	16	35
1	B	653/653 (100%)	592 (91%)	61 (9%)	13	29
All	All	1306/1306 (100%)	1191 (91%)	115 (9%)	14	32

All (115) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	SER
1	A	40	ARG
1	A	42	THR
1	A	49	LEU
1	A	60	LEU
1	A	82	GLU
1	A	87	SER

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Mol	Chain	Res	Type
1	A	88	VAL
1	A	107	ILE
1	A	120	TYR
1	A	180	LEU
1	A	202	VAL
1	A	211	TYR
1	A	223	LEU
1	A	230	ASP
1	A	246	LEU
1	A	273	THR
1	A	277	SER
1	A	278	SER
1	A	294	LEU
1	A	316	LEU
1	A	326	ASP
1	A	343	ARG
1	A	350	THR
1	A	377	ASN
1	A	385	CYS
1	A	399	LYS
1	A	410	LEU
1	A	419	SER
1	A	420	ASN
1	A	431	LEU
1	A	435	GLN
1	A	452	GLU
1	A	460	SER
1	A	471	ARG
1	A	472	CYS
1	A	473	SER
1	A	479	LEU
1	A	482	LEU
1	A	500	LEU
1	A	502	LYS
1	A	513	LYS
1	A	543	LEU
1	A	544	LEU
1	A	621	ASN
1	A	630	SER
1	A	663	ASP
1	A	669	ARG
1	A	679	ASN

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Mol	Chain	Res	Type
1	A	684	ARG
1	A	710	ASN
1	A	715	GLN
1	A	736	THR
1	A	745	SER
1	B	41	LYS
1	B	42	THR
1	B	49	LEU
1	B	60	LEU
1	B	63	ILE
1	B	71	LYS
1	B	88	VAL
1	B	90	LEU
1	B	91	GLU
1	B	94	THR
1	B	129	THR
1	B	143	ILE
1	B	156	THR
1	B	202	VAL
1	B	214	LEU
1	B	217	SER
1	B	223	LEU
1	B	230	ASP
1	B	236	ILE
1	B	246	LEU
1	B	263	ASN
1	B	273	THR
1	B	288	THR
1	B	295	ILE
1	B	300	LEU
1	B	302	ASP
1	B	314	GLN
1	B	326	ASP
1	B	332	GLU
1	B	334	SER
1	B	350	THR
1	B	377	ASN
1	B	385	CYS
1	B	389	ILE
1	B	392	LYS
1	B	397	ILE
1	B	420	ASN

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Mol	Chain	Res	Type
1	B	452	GLU
1	B	479	LEU
1	B	486	VAL
1	B	500	LEU
1	B	507	VAL
1	B	514	LEU
1	B	558	VAL
1	B	562	ASN
1	B	566	TYR
1	B	594	ILE
1	B	600	THR
1	B	619	VAL
1	B	620	ASP
1	B	621	ASN
1	B	630	SER
1	B	644	SER
1	B	677	GLU
1	B	679	ASN
1	B	680	LEU
1	B	684	ARG
1	B	710	ASN
1	B	715	GLN
1	B	726	VAL
1	B	733	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	GLN
1	A	169	ASN
1	A	247	GLN
1	A	263	ASN
1	A	377	ASN
1	A	420	ASN
1	A	483	HIS
1	A	508	GLN
1	A	533	HIS
1	A	592	HIS
1	A	621	ASN
1	A	679	ASN
1	A	710	ASN
1	A	715	GLN

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Mol	Chain	Res	Type
1	A	750	HIS
1	B	126	HIS
1	B	169	ASN
1	B	263	ASN
1	B	286	GLN
1	B	298	HIS
1	B	377	ASN
1	B	383	HIS
1	B	420	ASN
1	B	435	GLN
1	B	483	HIS
1	B	562	ASN
1	B	621	ASN
1	B	679	ASN
1	B	682	HIS
1	B	710	ASN
1	B	715	GLN
1	B	750	HIS
1	B	757	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	A	793	1	12,14,15	0.81	1 (8%)	15,19,21	2.08	4 (26%)
2	NAG	A	794	1	12,14,15	0.92	1 (8%)	15,19,21	2.02	6 (40%)
2	NAG	A	795	1	12,14,15	0.65	0	15,19,21	1.62	3 (20%)
2	NAG	A	796	1	12,14,15	0.84	1 (8%)	15,19,21	1.93	5 (33%)
3	B2Q	A	900	-	25,25,25	0.86	1 (4%)	35,35,35	2.07	10 (28%)
2	NAG	B	793	1	12,14,15	0.73	0	15,19,21	1.42	5 (33%)
2	NAG	B	794	1	12,14,15	0.80	0	15,19,21	1.42	2 (13%)
2	NAG	B	796	1	12,14,15	0.84	1 (8%)	15,19,21	1.38	2 (13%)
2	NAG	B	797	1	12,14,15	0.60	0	15,19,21	1.52	3 (20%)
3	B2Q	B	900	-	25,25,25	1.39	2 (8%)	35,35,35	2.91	12 (34%)

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
2	NAG	A	793	1	-	0/6/23/26	0/1/1/1
2	NAG	A	794	1	-	0/6/23/26	0/1/1/1
2	NAG	A	795	1	-	0/6/23/26	0/1/1/1
2	NAG	A	796	1	-	0/6/23/26	0/1/1/1
3	B2Q	A	900	-	2/2/4/4	0/8/33/33	0/0/3/3
2	NAG	B	793	1	-	0/6/23/26	0/1/1/1
2	NAG	B	794	1	-	0/6/23/26	0/1/1/1
2	NAG	B	796	1	-	0/6/23/26	0/1/1/1
2	NAG	B	797	1	-	0/6/23/26	0/1/1/1
3	B2Q	B	900	-	2/2/4/4	0/8/33/33	0/0/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	900	B2Q	CL-CR	-5.39	1.47	1.53
3	A	900	B2Q	CT-CR	3.01	1.54	1.52
3	B	900	B2Q	CT-CR	3.00	1.54	1.52
2	B	796	NAG	O5-C5	-2.65	1.40	1.45
2	A	793	NAG	O5-C5	-2.57	1.40	1.45
2	A	794	NAG	O5-C5	-2.28	1.41	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	796	NAG	O5-C5	-2.06	1.41	1.45

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	900	B2Q	CK-CL-CR	7.12	119.06	110.26
3	B	900	B2Q	CM-CL-CK	-6.95	103.60	111.67
3	B	900	B2Q	CL-CK-NJ	-6.67	103.22	112.13
3	B	900	B2Q	CN-CM-CL	6.22	140.61	115.47
3	B	900	B2Q	CT-CR-NS	-5.07	95.39	111.00
3	A	900	B2Q	CT-CR-NS	-5.07	95.42	111.00
2	A	795	NAG	C3-C2-N2	4.51	118.63	111.76
2	A	793	NAG	O5-C5-C6	-4.50	102.26	106.98
3	A	900	B2Q	CL-CK-NJ	-3.99	106.80	112.13
2	A	796	NAG	C3-C4-C5	-3.97	103.11	110.20
2	A	794	NAG	O5-C5-C6	3.92	111.10	106.98
3	B	900	B2Q	CF-CD-CE	-3.84	118.01	121.63
3	A	900	B2Q	CF-CD-CE	-3.83	118.01	121.63
3	A	900	B2Q	CM-CL-CR	-3.74	107.01	112.07
3	B	900	B2Q	CH-NJ-CE	3.61	118.55	111.08
3	A	900	B2Q	CH-NJ-CE	3.60	118.54	111.08
2	B	797	NAG	O5-C5-C6	3.57	110.72	106.98
3	A	900	B2Q	CL-CR-NS	-3.39	100.59	111.39
2	A	793	NAG	C3-C2-N2	3.29	116.76	111.76
3	A	900	B2Q	CN-CM-CL	3.21	128.44	115.47
2	A	793	NAG	O5-C5-C4	3.16	114.66	110.65
2	A	796	NAG	C3-C2-N2	-3.08	107.08	111.76
3	B	900	B2Q	CL-CR-NS	-3.06	101.65	111.39
2	B	796	NAG	O5-C5-C6	-2.98	103.85	106.98
2	B	797	NAG	C3-C4-C5	-2.97	104.90	110.20
2	B	794	NAG	C3-C2-N2	-2.91	107.34	111.76
2	B	796	NAG	C3-C2-N2	-2.88	107.37	111.76
2	A	794	NAG	C3-C2-N2	-2.82	107.46	111.76
2	A	796	NAG	O3-C3-C2	-2.68	103.46	109.09
2	A	795	NAG	C4-C3-C2	-2.63	104.88	111.32
2	A	794	NAG	C6-C5-C4	-2.57	106.79	113.00
3	B	900	B2Q	CM-CL-CR	2.56	115.54	112.07
2	A	794	NAG	C3-C4-C5	-2.52	105.70	110.20
2	B	794	NAG	O5-C5-C6	2.50	109.60	106.98
3	B	900	B2Q	OB-CB-CV	2.44	118.96	115.42
3	B	900	B2Q	CT-CR-CL	-2.42	106.77	111.33
3	A	900	B2Q	OB-CB-CV	2.42	118.93	115.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	793	NAG	C3-C4-C5	2.41	114.50	110.20
3	A	900	B2Q	OW-CV-CB	2.40	118.91	115.42
3	B	900	B2Q	OW-CV-CB	2.36	118.84	115.42
2	A	794	NAG	C4-C3-C2	-2.34	105.58	111.32
2	B	793	NAG	O3-C3-C2	2.31	113.95	109.09
3	A	900	B2Q	CK-CL-CR	2.29	113.09	110.26
2	B	793	NAG	C4-C3-C2	-2.26	105.79	111.32
2	A	796	NAG	O7-C7-N2	2.26	126.62	121.90
2	A	796	NAG	O4-C4-C5	2.25	115.20	109.28
2	A	795	NAG	O5-C5-C6	2.15	109.24	106.98
2	B	793	NAG	O5-C5-C4	2.15	113.38	110.65
2	B	797	NAG	C2-N2-C7	-2.13	119.51	123.09
2	B	793	NAG	C6-C5-C4	-2.13	107.86	113.00
2	A	794	NAG	O7-C7-C8	-2.03	118.08	122.04
2	B	793	NAG	C8-C7-N2	2.01	120.03	116.11

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	900	B2Q	CE
3	B	900	B2Q	CL
3	A	900	B2Q	CE
3	A	900	B2Q	CL

There are no torsion outliers.

There are no ring outliers.

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, 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	728/728 (100%)	0.35	11 (1%) 70 75	27, 44, 68, 93	0
1	B	728/728 (100%)	0.25	11 (1%) 70 75	27, 44, 65, 84	0
All	All	1456/1456 (100%)	0.30	22 (1%) 70 75	27, 44, 67, 93	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	140	ARG	4.4
1	A	140	ARG	3.6
1	B	97	GLU	3.4
1	A	141	GLN	3.4
1	A	296	GLY	3.3
1	A	145	GLU	3.2
1	A	74	ASN	3.1
1	A	99	GLY	3.0
1	B	279	VAL	2.8
1	A	142	LEU	2.6
1	A	471	ARG	2.6
1	A	615	LYS	2.5
1	B	145	GLU	2.5
1	B	74	ASN	2.5
1	B	278	SER	2.2
1	B	144	THR	2.2
1	B	597	ARG	2.2
1	B	189	GLY	2.1
1	A	333	SER	2.1
1	A	180	LEU	2.1
1	B	392	LYS	2.1
1	B	340	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	B	797	14/15	0.32	3.58	81,83,85,85	0
2	NAG	B	793	14/15	0.33	3.11	72,79,81,82	0
3	B2Q	B	900	23/23	0.22	1.84	41,48,49,49	0
2	NAG	B	796	14/15	0.26	1.56	57,62,63,65	0
2	NAG	A	795	14/15	0.27	1.12	83,85,87,87	0
2	NAG	A	794	14/15	0.24	0.74	46,48,53,54	0
3	B2Q	A	900	23/23	0.19	0.11	42,45,46,47	0
2	NAG	B	794	14/15	0.18	0.06	53,55,56,57	0
2	NAG	A	793	14/15	0.19	-0.47	71,73,76,76	0
2	NAG	A	796	14/15	0.18	-0.50	49,52,54,56	0

6.5 Other polymers ⓘ

There are no such residues in this entry.