



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:47 AM UTC

PDB ID : 6NA6 / pdb_00006na6
Title : Serial Femtosecond X-ray Crystallography Structure of ECR in complex with NADPH
Authors : DeMirici, H.
Deposited on : 2018-12-05
Resolution : 2.10 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

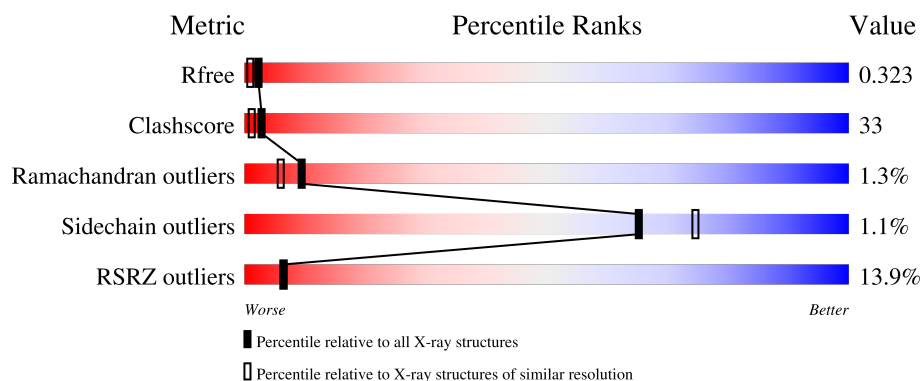
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	445	16% 53% 43% ..
1	B	445	10% 57% 42% .
1	C	445	11% 56% 42% .
1	D	445	19% 59% 40% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NDP	A	501	X	-	-	-
2	NDP	B	501	X	-	-	-
2	NDP	C	501	X	-	-	-
2	NDP	D	501	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14569 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.

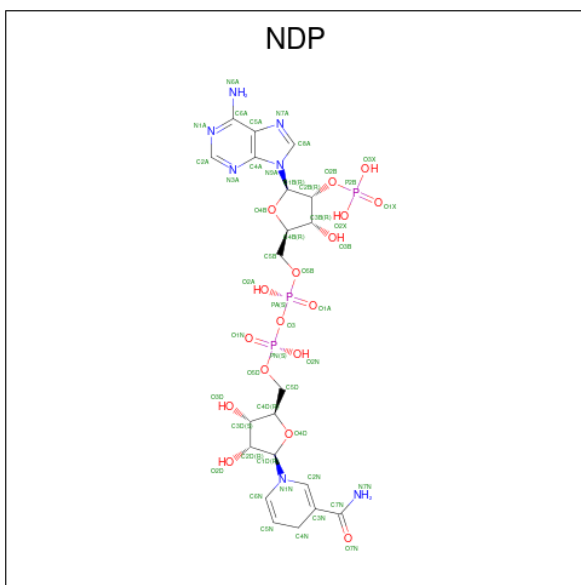
- Molecule 1 is a protein called Putative crotonyl-CoA reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	1	0
			3443	2166	620	644	13			
1	B	445	Total	C	N	O	S	6	2	0
			3449	2169	621	646	13			
1	D	445	Total	C	N	O	S	6	2	0
			3449	2169	621	646	13			
1	C	445	Total	C	N	O	S	6	2	0
			3449	2169	621	646	13			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ARG	-	expression tag	UNP E4N096
A	0	HIS	-	expression tag	UNP E4N096
B	-1	ARG	-	expression tag	UNP E4N096
B	0	HIS	-	expression tag	UNP E4N096
D	-1	ARG	-	expression tag	UNP E4N096
D	0	HIS	-	expression tag	UNP E4N096
C	-1	ARG	-	expression tag	UNP E4N096
C	0	HIS	-	expression tag	UNP E4N096

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	D	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	C	1	Total 48	C 21	N 7	O 17	P 3	0	0

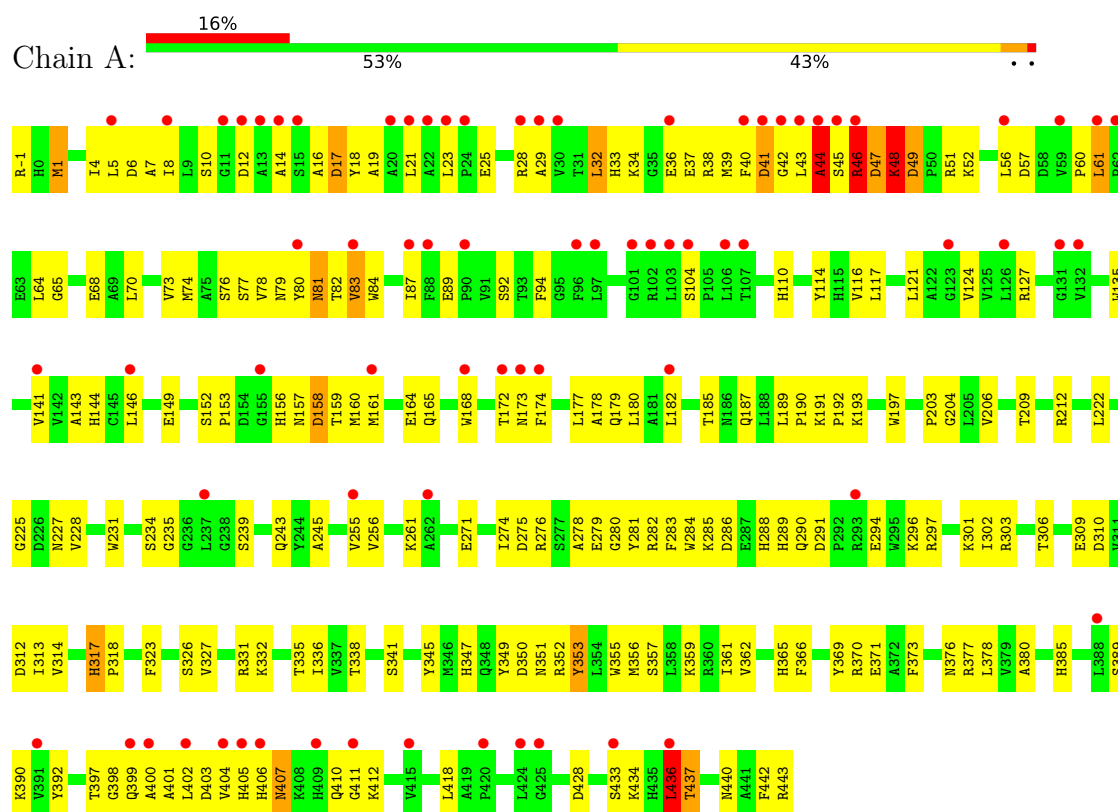
- Molecule 3 is water.

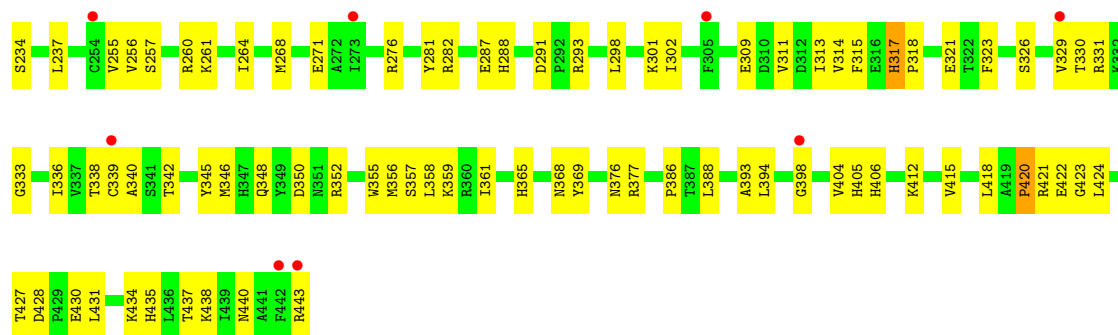
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	119	Total O 119 119	0	0
3	B	152	Total O 152 152	0	0
3	D	145	Total O 145 145	0	0
3	C	171	Total O 171 171	0	0

3 Residue-property plots

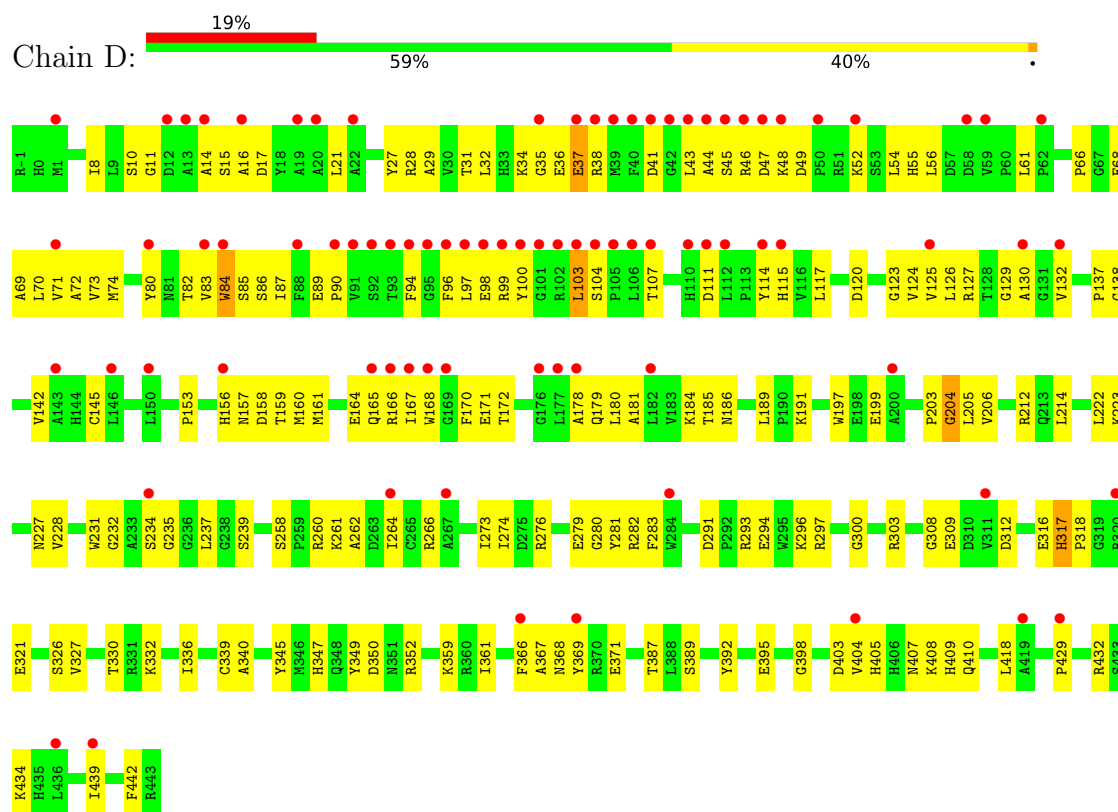
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: Putative crotonyl-CoA reductase

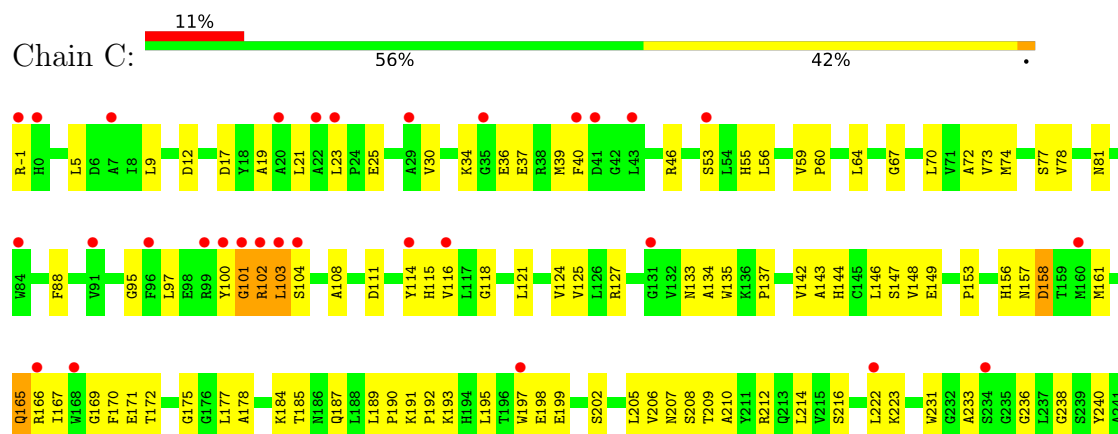


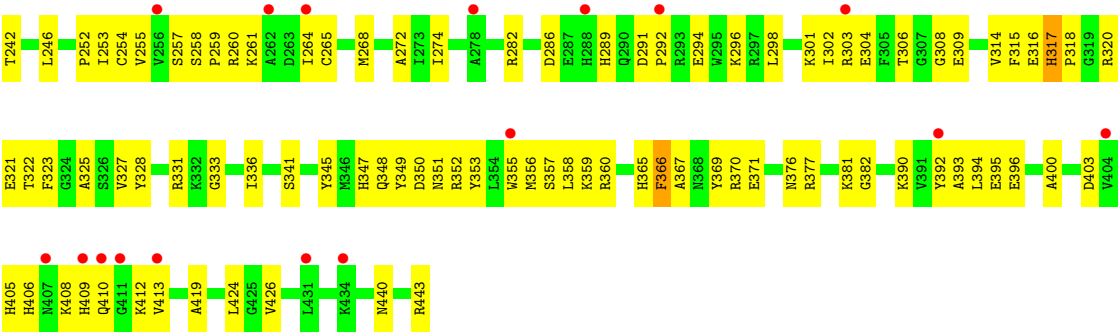


● Molecule 1: Putative crotonyl-CoA reductase



● Molecule 1: Putative crotonyl-CoA reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	109.76Å 78.05Å 138.00Å 90.00° 107.76° 90.00°	Depositor
Resolution (Å)	31.82 – 2.10 31.82 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.6 (31.82-2.10) 84.0 (31.82-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.19 (at 1.62Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.274 , 0.333 0.264 , 0.323	Depositor DCC
R_{free} test set	1851 reflections (1.14%)	wwPDB-VP
Wilson B-factor (Å ²)	13.5	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 74.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	14569	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 45.68 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2758e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

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

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	0.77	5/3525 (0.1%)	0.79	8/4779 (0.2%)
1	B	0.41	0/3531	0.63	0/4787
1	C	0.42	0/3531	0.66	0/4787
1	D	0.39	0/3531	0.62	0/4787
All	All	0.52	5/14118 (0.0%)	0.68	8/19140 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	D	0	1
All	All	0	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	353	TYR	C-O	-5.37	1.17	1.24
1	A	61	LEU	C-O	-5.31	1.16	1.24
1	A	1	MET	C-O	-5.21	1.18	1.24
1	A	428	ASP	N-CA	-5.14	1.41	1.46
1	A	83	VAL	C-O	-5.11	1.18	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	ASP	N-CA-C	7.19	119.98	111.71
1	A	1	MET	N-CA-C	-6.47	104.22	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	17	ASP	N-CA-C	-6.23	104.41	111.14
1	A	48	LYS	N-CA-C	-5.84	101.20	109.96
1	A	46	ARG	N-CA-C	5.59	120.97	114.04
1	A	436	LEU	CA-C-N	5.56	129.16	120.82
1	A	436	LEU	C-N-CA	5.56	129.16	120.82
1	A	44	ALA	N-CA-C	5.03	121.51	110.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	338	THR	Peptide
1	D	84	TRP	Peptide

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	3443	0	3375	264	0
1	B	3449	0	3379	210	0
1	C	3449	0	3379	245	0
1	D	3449	0	3379	228	0
2	A	48	0	25	11	0
2	B	48	0	25	3	0
2	C	48	0	24	5	0
2	D	48	0	25	16	0
3	A	119	0	0	89	0
3	B	152	0	0	71	0
3	C	171	0	0	92	0
3	D	145	0	0	86	0
All	All	14569	0	13611	903	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 33.

All (903) 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:142:VAL:HG22	3:C:606:HOH:O	1.29	1.30
1:B:115:HIS:HE1	3:B:608:HOH:O	1.11	1.27
1:D:71:VAL:HG12	3:D:605:HOH:O	1.25	1.26
1:A:190:PRO:HA	3:A:605:HOH:O	1.26	1.24
1:B:415:VAL:HG13	3:B:607:HOH:O	1.36	1.23
2:A:501:NDP:O4B	2:A:501:NDP:C1B	1.63	1.22
1:A:127:ARG:CA	3:A:607:HOH:O	1.86	1.22
1:A:282:ARG:HG2	3:A:603:HOH:O	1.40	1.22
1:A:127:ARG:CB	3:A:607:HOH:O	1.88	1.20
1:A:127:ARG:HB2	3:A:607:HOH:O	1.42	1.19
1:D:43:LEU:C	3:D:601:HOH:O	1.82	1.19
1:B:115:HIS:CE1	3:B:608:HOH:O	1.87	1.16
1:C:147:SER:HB2	3:C:609:HOH:O	1.46	1.16
1:A:418:LEU:HG	3:A:613:HOH:O	1.45	1.14
1:C:147:SER:CB	3:C:609:HOH:O	1.95	1.14
1:C:331:ARG:NE	3:C:601:HOH:O	1.81	1.14
1:C:331:ARG:NH2	3:C:601:HOH:O	1.81	1.13
1:A:10:SER:HB3	3:A:610:HOH:O	1.47	1.13
1:C:209:THR:N	3:C:603:HOH:O	1.81	1.13
1:A:261:LYS:NZ	3:A:601:HOH:O	1.80	1.12
1:B:412:LYS:NZ	3:B:601:HOH:O	1.80	1.12
1:D:303:ARG:HG2	3:D:618:HOH:O	1.47	1.12
1:B:226:ASP:N	3:B:602:HOH:O	1.81	1.11
1:B:393:ALA:C	3:B:607:HOH:O	1.90	1.11
1:D:44:ALA:N	3:D:601:HOH:O	1.82	1.11
1:B:394:LEU:N	3:B:607:HOH:O	1.84	1.11
1:D:15:SER:N	3:D:604:HOH:O	1.84	1.10
1:D:43:LEU:CB	3:D:601:HOH:O	1.99	1.10
1:D:100:TYR:N	3:D:608:HOH:O	1.86	1.07
1:C:254:CYS:C	3:C:605:HOH:O	1.98	1.07
1:A:42:GLY:O	1:A:43:LEU:HD23	1.54	1.07
1:D:115:HIS:NE2	3:D:612:HOH:O	1.88	1.06
1:A:127:ARG:N	3:A:607:HOH:O	1.83	1.06
1:D:189:LEU:C	3:D:613:HOH:O	1.97	1.06
1:C:331:ARG:CZ	3:C:601:HOH:O	2.03	1.05
1:A:418:LEU:CG	3:A:613:HOH:O	1.99	1.05
1:A:418:LEU:CA	3:A:613:HOH:O	2.03	1.04
1:B:340:ALA:HB1	2:B:501:NDP:H6N	1.39	1.04
1:A:418:LEU:N	3:A:613:HOH:O	1.89	1.04
1:C:142:VAL:N	3:C:606:HOH:O	1.87	1.03
1:D:186:ASN:ND2	3:D:606:HOH:O	1.85	1.02
1:D:14:ALA:C	3:D:604:HOH:O	2.02	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:283:PHE:N	3:D:614:HOH:O	1.93	1.01
1:C:351:ASN:ND2	1:C:355:TRP:HE3	1.59	1.00
1:A:282:ARG:NH2	3:A:616:HOH:O	1.93	0.99
1:B:222:LEU:HA	3:B:661:HOH:O	1.62	0.99
1:B:281:TYR:OH	3:B:603:HOH:O	1.81	0.98
1:A:64:LEU:O	3:A:602:HOH:O	1.81	0.98
1:A:437:THR:OG1	3:A:604:HOH:O	1.82	0.97
1:B:394:LEU:CA	3:B:607:HOH:O	2.12	0.97
1:D:153:PRO:HG2	3:D:725:HOH:O	1.64	0.97
1:A:281:TYR:O	3:A:603:HOH:O	1.82	0.96
1:A:373:PHE:CE1	3:D:725:HOH:O	2.15	0.96
1:B:186:ASN:OD1	3:B:605:HOH:O	1.82	0.96
1:D:403:ASP:OD2	3:D:603:HOH:O	1.82	0.96
1:C:413:VAL:O	3:C:604:HOH:O	1.82	0.96
1:A:281:TYR:O	3:A:608:HOH:O	1.83	0.96
1:B:170:PHE:O	3:B:608:HOH:O	1.84	0.96
1:B:282:ARG:NH1	3:B:616:HOH:O	1.98	0.95
1:D:181:ALA:HA	3:D:633:HOH:O	1.63	0.95
1:D:260:ARG:NH1	3:D:607:HOH:O	1.86	0.95
1:C:207:ASN:C	3:C:603:HOH:O	2.08	0.95
1:A:135:TRP:NE1	3:A:609:HOH:O	1.83	0.95
1:A:6:ASP:O	3:A:610:HOH:O	1.84	0.95
1:A:7:ALA:C	3:A:610:HOH:O	2.07	0.95
1:A:189:LEU:O	3:A:605:HOH:O	1.82	0.95
1:B:443:ARG:O	3:B:606:HOH:O	1.83	0.95
1:D:312:ASP:OD1	3:D:602:HOH:O	1.82	0.95
1:D:72:ALA:O	3:D:605:HOH:O	1.85	0.95
1:A:7:ALA:O	3:A:610:HOH:O	1.83	0.95
1:A:48:LYS:HB3	1:A:87:ILE:HD11	1.46	0.95
1:C:320:ARG:NE	3:C:613:HOH:O	1.95	0.94
1:D:308:GLY:C	3:D:618:HOH:O	2.09	0.94
1:B:10:SER:OG	3:B:609:HOH:O	1.85	0.94
1:B:25:GLU:HG3	3:B:627:HOH:O	1.66	0.94
1:C:254:CYS:N	3:C:605:HOH:O	1.99	0.94
1:C:351:ASN:ND2	1:C:355:TRP:CE3	2.35	0.94
1:A:92:SER:N	3:A:606:HOH:O	1.83	0.94
1:C:292:PRO:HD2	3:C:617:HOH:O	1.67	0.93
1:D:389:SER:CB	3:D:622:HOH:O	2.14	0.93
1:C:184:LYS:HE3	3:C:609:HOH:O	1.69	0.93
1:B:333:GLY:HA3	1:D:161:MET:HE1	1.50	0.93
1:B:-1:ARG:HB3	3:B:619:HOH:O	1.68	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:LEU:O	3:B:611:HOH:O	1.88	0.92
1:B:394:LEU:HA	3:B:607:HOH:O	1.69	0.91
1:D:138:GLY:O	3:D:610:HOH:O	1.87	0.91
1:A:341:SER:OG	1:C:355:TRP:HZ3	1.54	0.91
1:C:210:ALA:N	3:C:603:HOH:O	2.03	0.91
1:A:135:TRP:C	3:A:619:HOH:O	2.13	0.91
1:C:252:PRO:O	3:C:608:HOH:O	1.89	0.91
1:A:64:LEU:HD12	3:A:602:HOH:O	1.70	0.91
1:D:14:ALA:CB	3:D:604:HOH:O	2.19	0.90
1:A:312:ASP:OD2	3:A:612:HOH:O	1.89	0.89
1:A:433:SER:C	3:A:620:HOH:O	2.14	0.89
1:A:225:GLY:O	1:B:377:ARG:NH2	2.05	0.89
1:D:409:HIS:O	3:D:611:HOH:O	1.88	0.89
1:A:46:ARG:HB3	1:A:406:HIS:NE2	1.88	0.89
1:B:437:THR:OG1	3:B:610:HOH:O	1.85	0.89
1:C:440:ASN:O	3:C:607:HOH:O	1.88	0.89
1:B:79:ASN:OD1	3:B:612:HOH:O	1.92	0.88
1:C:254:CYS:O	3:C:605:HOH:O	1.85	0.87
1:A:271:GLU:OE2	3:A:614:HOH:O	1.90	0.87
1:B:10:SER:CB	3:B:609:HOH:O	2.22	0.87
1:A:314:VAL:HG11	1:A:326:SER:HB2	1.56	0.87
1:C:-1:ARG:NH1	3:C:622:HOH:O	2.06	0.87
1:D:172:THR:C	3:D:612:HOH:O	2.16	0.87
1:B:330:THR:HG21	1:B:336:ILE:HD11	1.56	0.87
1:D:165:GLN:NE2	3:D:617:HOH:O	1.96	0.87
1:C:254:CYS:O	3:C:610:HOH:O	1.91	0.87
1:D:189:LEU:O	3:D:613:HOH:O	1.88	0.86
1:B:430:GLU:OE2	3:B:614:HOH:O	1.93	0.86
1:D:164:GLU:OE2	3:D:615:HOH:O	1.93	0.86
1:C:206:VAL:O	3:C:603:HOH:O	1.94	0.86
2:A:501:NDP:O1N	3:A:617:HOH:O	1.93	0.86
1:B:-1:ARG:CB	3:B:619:HOH:O	2.22	0.86
1:D:99:ARG:C	3:D:608:HOH:O	2.15	0.86
1:D:43:LEU:HB2	3:D:601:HOH:O	1.68	0.85
1:A:152:SER:O	3:A:615:HOH:O	1.92	0.85
1:A:261:LYS:HG2	3:A:601:HOH:O	1.75	0.85
1:D:16:ALA:N	3:D:604:HOH:O	2.09	0.85
1:A:81:ASN:N	3:A:623:HOH:O	2.10	0.85
1:D:392:TYR:OH	1:D:403:ASP:OD2	1.95	0.85
1:C:304:GLU:OE1	3:C:612:HOH:O	1.94	0.85
1:D:297:ARG:NH1	3:D:621:HOH:O	2.08	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:SER:HB2	1:C:355:TRP:CH2	2.12	0.84
1:A:190:PRO:CA	3:A:605:HOH:O	1.93	0.84
1:B:199:GLU:OE2	3:B:615:HOH:O	1.94	0.84
1:A:271:GLU:CD	3:A:614:HOH:O	2.19	0.84
1:D:142:VAL:HG22	3:D:613:HOH:O	1.76	0.83
1:D:172:THR:O	3:D:612:HOH:O	1.95	0.83
1:D:191:LYS:NZ	3:D:609:HOH:O	1.87	0.83
1:A:135:TRP:O	3:A:619:HOH:O	1.97	0.83
1:B:313:ILE:HD11	3:B:661:HOH:O	1.79	0.83
1:C:255:VAL:HA	3:C:610:HOH:O	1.78	0.83
1:A:282:ARG:CG	3:A:603:HOH:O	2.07	0.82
1:A:341:SER:OG	1:C:355:TRP:CZ3	2.32	0.82
1:A:373:PHE:CZ	3:D:725:HOH:O	2.30	0.81
1:B:167:ILE:HG13	1:B:170:PHE:HB3	1.61	0.81
1:D:99:ARG:N	3:D:608:HOH:O	2.14	0.81
1:A:433:SER:O	3:A:620:HOH:O	1.98	0.81
1:D:434:LYS:NZ	3:D:619:HOH:O	1.99	0.80
1:C:258:SER:HB3	1:C:261:LYS:HD3	1.63	0.80
1:D:71:VAL:CG1	3:D:605:HOH:O	1.99	0.80
1:C:72:ALA:O	3:C:615:HOH:O	2.00	0.80
1:C:393:ALA:HB3	3:C:620:HOH:O	1.81	0.80
1:B:133:ASN:HB2	3:B:707:HOH:O	1.82	0.80
1:B:149:GLU:OE1	3:B:617:HOH:O	2.00	0.79
1:C:409:HIS:O	3:C:614:HOH:O	2.00	0.79
1:C:426:VAL:O	3:C:616:HOH:O	2.01	0.79
1:B:61:LEU:O	3:B:618:HOH:O	2.00	0.79
1:A:80:TYR:OH	2:A:501:NDP:O2A	1.99	0.79
1:A:48:LYS:CB	1:A:87:ILE:HD11	2.13	0.79
1:B:234:SER:OG	3:B:613:HOH:O	1.92	0.79
1:A:282:ARG:HA	3:A:603:HOH:O	1.82	0.78
1:C:301:LYS:CD	3:C:612:HOH:O	2.31	0.78
1:B:-1:ARG:CG	3:B:619:HOH:O	2.31	0.78
1:D:327:VAL:HG21	1:D:349:TYR:HB2	1.66	0.78
1:C:291:ASP:OD2	3:C:617:HOH:O	2.01	0.78
1:D:234:SER:HB2	1:D:261:LYS:HD2	1.66	0.78
1:C:114:TYR:OH	3:C:618:HOH:O	2.02	0.77
1:C:208:SER:C	3:C:603:HOH:O	2.17	0.77
1:D:14:ALA:HB1	3:D:604:HOH:O	1.83	0.77
1:A:81:ASN:OD1	3:A:623:HOH:O	2.01	0.76
1:B:281:TYR:O	3:B:616:HOH:O	2.02	0.76
1:C:301:LYS:HE2	3:C:612:HOH:O	1.83	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:THR:O	3:A:624:HOH:O	2.02	0.76
1:D:186:ASN:CG	3:D:606:HOH:O	2.21	0.76
1:B:443:ARG:C	3:B:606:HOH:O	2.28	0.76
1:D:166:ARG:HG2	3:D:634:HOH:O	1.84	0.76
1:D:308:GLY:O	3:D:618:HOH:O	1.99	0.76
1:A:10:SER:CB	3:A:610:HOH:O	2.15	0.76
1:C:202:SER:O	3:C:619:HOH:O	2.04	0.75
1:D:158:ASP:N	3:D:627:HOH:O	2.17	0.75
1:D:234:SER:CB	1:D:261:LYS:HD2	2.15	0.75
1:C:253:ILE:C	3:C:605:HOH:O	2.27	0.75
1:C:254:CYS:C	3:C:610:HOH:O	2.27	0.75
1:A:373:PHE:CD1	3:D:725:HOH:O	2.39	0.75
1:A:33:HIS:HB2	1:A:36:GLU:OE2	1.87	0.75
1:C:19:ALA:O	1:C:127:ARG:NH2	2.19	0.75
1:C:393:ALA:O	3:C:620:HOH:O	2.04	0.75
1:C:291:ASP:HA	3:C:617:HOH:O	1.87	0.74
1:B:63:GLU:HG3	1:B:109:ARG:HD3	1.68	0.74
1:C:255:VAL:CA	3:C:610:HOH:O	2.35	0.74
1:A:436:LEU:HB3	3:A:620:HOH:O	1.88	0.74
1:D:32:LEU:HD21	1:D:86:SER:HA	1.70	0.74
1:D:115:HIS:CD2	3:D:612:HOH:O	2.32	0.74
1:A:338:THR:HG21	1:C:351:ASN:HD21	1.50	0.73
1:B:10:SER:HB3	3:B:609:HOH:O	1.83	0.73
1:A:84:TRP:HA	1:A:87:ILE:HG22	1.70	0.73
1:C:396:GLU:HB2	3:C:620:HOH:O	1.86	0.73
1:D:279:GLU:N	3:D:630:HOH:O	2.21	0.73
1:B:231:TRP:HA	1:B:255:VAL:HG12	1.70	0.73
1:D:389:SER:HB2	3:D:622:HOH:O	1.82	0.73
1:C:345:TYR:OH	3:C:621:HOH:O	2.05	0.73
1:D:300:GLY:HA2	1:D:303:ARG:HH21	1.54	0.72
1:B:404:VAL:O	3:B:620:HOH:O	2.08	0.72
1:D:186:ASN:OD1	3:D:606:HOH:O	2.06	0.72
1:C:169:GLY:H	1:C:172:THR:HG22	1.54	0.72
1:C:393:ALA:CA	3:C:620:HOH:O	2.36	0.72
1:B:257:SER:CB	3:B:604:HOH:O	2.29	0.72
1:D:274:ILE:HD11	1:D:281:TYR:HE2	1.54	0.72
1:A:149:GLU:HB2	1:A:164:GLU:HG3	1.70	0.72
1:C:395:GLU:O	3:C:623:HOH:O	2.08	0.72
1:A:46:ARG:CB	1:A:406:HIS:NE2	2.52	0.72
1:A:81:ASN:N	3:A:625:HOH:O	2.07	0.72
1:A:350:ASP:OD1	1:A:352:ARG:HG3	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LYS:HB3	1:A:87:ILE:CD1	2.20	0.71
1:A:157:ASN:OD1	1:D:368:ASN:ND2	2.21	0.71
1:B:229:LEU:H	1:B:311:VAL:HG21	1.55	0.71
1:D:317:HIS:HB2	2:D:501:NDP:H3D	1.71	0.71
1:A:46:ARG:HB3	1:A:406:HIS:CD2	2.25	0.71
1:B:264:ILE:HD11	3:B:624:HOH:O	1.89	0.71
1:A:19:ALA:O	1:A:127:ARG:NH2	2.24	0.70
1:C:382:GLY:HA2	3:C:702:HOH:O	1.91	0.70
1:B:164:GLU:HG3	1:B:166:ARG:HH12	1.56	0.70
1:A:168:TRP:HA	1:A:172:THR:HG22	1.73	0.70
1:B:99:ARG:HB2	3:B:655:HOH:O	1.92	0.70
1:B:66:PRO:HD2	1:B:106:LEU:HD11	1.73	0.70
1:B:231:TRP:CE3	1:B:255:VAL:HG11	2.26	0.70
1:D:189:LEU:CB	3:D:613:HOH:O	2.39	0.70
1:A:127:ARG:C	3:A:607:HOH:O	2.24	0.70
1:A:7:ALA:CA	3:A:610:HOH:O	2.36	0.70
1:D:258:SER:OG	1:D:261:LYS:NZ	2.25	0.69
1:C:207:ASN:O	3:C:603:HOH:O	2.04	0.69
1:A:70:LEU:HD11	1:A:180:LEU:HD13	1.73	0.69
1:C:206:VAL:HG11	1:C:412:LYS:HD3	1.73	0.69
1:A:65:GLY:HA2	3:A:602:HOH:O	1.93	0.68
1:A:228:VAL:HG12	1:A:313:ILE:HB	1.74	0.68
1:C:424:LEU:HB2	3:C:687:HOH:O	1.91	0.68
1:D:340:ALA:HB1	3:D:674:HOH:O	1.93	0.68
1:A:341:SER:CB	1:C:355:TRP:CZ3	2.76	0.68
1:A:18:TYR:HA	1:A:21:LEU:HD12	1.76	0.68
1:A:282:ARG:CA	3:A:603:HOH:O	2.41	0.68
1:D:115:HIS:NE2	1:D:172:THR:O	2.23	0.68
1:D:274:ILE:HD11	1:D:281:TYR:CE2	2.29	0.68
1:C:292:PRO:CD	3:C:617:HOH:O	2.33	0.68
1:D:189:LEU:HB2	3:D:613:HOH:O	1.94	0.68
1:C:124:VAL:O	3:C:615:HOH:O	2.12	0.68
1:C:406:HIS:HB2	1:C:408:LYS:HG3	1.75	0.67
1:B:98:GLU:O	1:B:99:ARG:NH2	2.25	0.67
1:D:32:LEU:HG	1:D:117:LEU:HD11	1.77	0.67
1:D:89:GLU:HG2	1:D:90:PRO:HA	1.75	0.67
1:A:51:ARG:NH1	1:A:402:LEU:HD13	2.09	0.67
1:D:309:GLU:OE2	1:C:377:ARG:NH2	2.24	0.67
1:C:274:ILE:HB	3:C:610:HOH:O	1.94	0.67
1:B:301:LYS:HB3	3:B:603:HOH:O	1.94	0.67
1:D:199:GLU:OE2	1:D:432:ARG:NH1	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:393:ALA:C	3:C:620:HOH:O	2.36	0.67
1:B:28:ARG:NE	1:B:56:LEU:CD2	2.58	0.67
1:D:98:GLU:HA	1:D:111:ASP:CG	2.20	0.67
1:C:317:HIS:H	1:C:318:PRO:CD	2.08	0.67
1:D:404:VAL:HA	1:D:409:HIS:CD2	2.29	0.66
1:C:392:TYR:OH	1:C:403:ASP:OD2	2.12	0.66
1:A:440:ASN:HB2	1:A:443:ARG:HG3	1.77	0.66
1:B:128:THR:OG1	1:B:132:VAL:HB	1.95	0.66
1:B:179:GLN:HB2	1:B:418:LEU:HG	1.78	0.66
1:A:357:SER:HB3	1:A:359:LYS:HE3	1.77	0.66
1:B:146:LEU:HD12	1:B:148:VAL:HG13	1.78	0.66
1:D:47:ASP:O	1:D:52:LYS:NZ	2.25	0.66
1:D:70:LEU:HG	1:D:126:LEU:HD12	1.77	0.66
1:C:240:TYR:OH	1:C:412:LYS:NZ	2.26	0.66
1:B:415:VAL:CG1	3:B:607:HOH:O	2.14	0.66
1:D:45:SER:HA	1:D:48:LYS:HE2	1.78	0.66
1:D:232:GLY:HA2	2:D:501:NDP:O1X	1.96	0.66
1:B:214:LEU:HD21	1:B:222:LEU:HD13	1.77	0.65
1:A:392:TYR:OH	1:A:403:ASP:OD1	2.05	0.65
1:A:51:ARG:HG3	1:A:398:GLY:HA3	1.79	0.65
1:B:423:GLY:O	3:B:621:HOH:O	2.15	0.65
1:C:198:GLU:O	3:C:625:HOH:O	2.15	0.65
1:A:127:ARG:O	3:A:607:HOH:O	2.12	0.65
1:B:146:LEU:CD1	1:B:148:VAL:HG13	2.26	0.65
1:B:330:THR:HG23	1:B:359:LYS:HD3	1.79	0.65
1:C:405:HIS:O	3:C:624:HOH:O	2.13	0.65
1:A:-1:ARG:HD3	3:A:643:HOH:O	1.96	0.65
1:B:331:ARG:HD3	3:B:626:HOH:O	1.96	0.65
1:D:389:SER:OG	3:D:622:HOH:O	2.11	0.65
1:A:336:ILE:HA	3:A:624:HOH:O	1.97	0.64
1:D:46:ARG:HA	3:D:628:HOH:O	1.97	0.64
1:C:393:ALA:N	3:C:620:HOH:O	2.30	0.64
1:A:38:ARG:C	1:A:40:PHE:H	2.05	0.64
1:A:48:LYS:CG	1:A:87:ILE:HD11	2.27	0.64
1:D:237:LEU:HD22	2:D:501:NDP:H1D	1.80	0.64
1:C:345:TYR:OH	3:C:611:HOH:O	1.92	0.64
1:A:135:TRP:N	3:A:619:HOH:O	2.29	0.64
1:D:84:TRP:HA	1:D:87:ILE:HB	1.78	0.64
1:A:281:TYR:C	3:A:608:HOH:O	2.33	0.64
1:B:-1:ARG:HG2	3:B:619:HOH:O	1.94	0.64
1:B:36:GLU:HG2	1:B:39:MET:SD	2.38	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:302:ILE:HD12	1:B:329:VAL:HG21	1.79	0.63
1:D:47:ASP:OD2	3:D:623:HOH:O	2.15	0.63
1:D:205:LEU:CD2	2:D:501:NDP:H5N	2.28	0.63
1:D:171:GLU:OE2	3:D:624:HOH:O	2.15	0.63
1:A:68:GLU:HB3	1:A:182:LEU:HD11	1.79	0.63
1:B:188:LEU:O	1:B:369:TYR:OH	2.14	0.63
1:B:54:LEU:HD12	1:B:398:GLY:HA2	1.80	0.63
2:A:501:NDP:O2A	3:A:627:HOH:O	2.16	0.63
1:D:167:ILE:HG22	1:D:171:GLU:OE1	1.97	0.63
1:C:301:LYS:HD2	3:C:612:HOH:O	1.96	0.63
1:A:261:LYS:CG	3:A:601:HOH:O	2.41	0.63
1:C:101:GLY:O	1:C:103:LEU:N	2.31	0.63
1:C:144:HIS:HA	3:C:677:HOH:O	1.98	0.62
1:B:213:GLN:HG2	1:B:315:PHE:CE2	2.34	0.62
1:A:271:GLU:OE1	3:A:614:HOH:O	2.16	0.62
1:B:154:ASP:HB3	1:D:332:LYS:HZ3	1.64	0.62
1:D:392:TYR:OH	3:D:603:HOH:O	2.06	0.62
1:A:79:ASN:OD1	1:A:412:LYS:NZ	2.23	0.62
1:D:124:VAL:N	3:D:605:HOH:O	2.28	0.62
1:C:146:LEU:HD11	1:C:148:VAL:HG13	1.80	0.62
1:C:64:LEU:HD11	1:C:70:LEU:HB2	1.82	0.62
1:C:214:LEU:HD13	1:C:315:PHE:HD2	1.64	0.61
1:D:227:ASN:HB2	3:D:602:HOH:O	1.99	0.61
1:A:341:SER:HB2	1:C:355:TRP:HH2	1.59	0.61
1:D:37:GLU:HG3	1:D:38:ARG:H	1.64	0.61
1:C:255:VAL:HG22	3:C:610:HOH:O	1.98	0.61
1:B:166:ARG:HB3	1:B:171:GLU:HB2	1.82	0.61
1:C:101:GLY:HA3	1:C:111:ASP:HB2	1.83	0.61
1:C:104:SER:CB	3:C:602:HOH:O	2.38	0.61
1:B:331:ARG:NH1	3:B:626:HOH:O	2.32	0.61
1:D:303:ARG:CG	3:D:618:HOH:O	2.23	0.61
1:D:206:VAL:HG22	3:D:641:HOH:O	2.00	0.60
1:B:355:TRP:CE3	1:D:340:ALA:HA	2.36	0.60
1:C:212:ARG:HA	1:C:216:SER:HB2	1.81	0.60
1:C:223:LYS:NZ	3:C:626:HOH:O	2.15	0.60
1:D:167:ILE:H	1:D:171:GLU:HB2	1.66	0.60
1:A:350:ASP:HB3	1:A:353:TYR:HD1	1.66	0.60
1:C:265:CYS:HA	1:C:268:MET:HE3	1.83	0.60
1:B:63:GLU:HG3	1:B:109:ARG:CD	2.31	0.60
1:D:184:LYS:HD3	3:D:606:HOH:O	2.01	0.60
1:A:39:MET:HE2	1:A:40:PHE:CZ	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:367:ALA:HB1	1:C:371:GLU:HB2	1.83	0.60
1:B:257:SER:HB2	3:B:604:HOH:O	1.97	0.60
1:D:214:LEU:HD21	1:D:222:LEU:HD13	1.83	0.60
1:A:341:SER:HB2	1:C:355:TRP:CZ3	2.36	0.60
1:D:199:GLU:CD	1:D:432:ARG:HH12	2.09	0.60
1:D:303:ARG:CD	3:D:618:HOH:O	2.49	0.59
1:C:351:ASN:HB3	1:C:355:TRP:CZ3	2.37	0.59
1:B:28:ARG:CZ	1:B:56:LEU:HD23	2.32	0.59
1:D:261:LYS:HE2	2:D:501:NDP:O2X	2.02	0.59
1:C:257:SER:HB2	2:C:501:NDP:O2X	2.02	0.59
1:D:100:TYR:HD1	1:D:103:LEU:HD13	1.66	0.59
1:C:102:ARG:HA	1:C:108:ALA:HB2	1.83	0.59
1:A:191:LYS:O	1:A:193:LYS:NZ	2.36	0.59
1:D:44:ALA:O	3:D:601:HOH:O	2.15	0.59
1:B:79:ASN:HA	3:B:612:HOH:O	2.02	0.59
1:A:16:ALA:CB	3:A:618:HOH:O	2.51	0.58
1:A:23:LEU:HD13	1:A:61:LEU:CD2	2.33	0.58
1:D:123:GLY:HA3	3:D:605:HOH:O	2.01	0.58
1:D:340:ALA:HB3	2:D:501:NDP:H2D	1.85	0.58
1:A:7:ALA:HA	3:A:610:HOH:O	2.01	0.58
1:B:195:LEU:C	3:B:611:HOH:O	2.42	0.58
1:D:38:ARG:HA	1:D:41:ASP:HB2	1.86	0.58
1:D:296:LYS:HE2	1:D:303:ARG:HH22	1.68	0.58
1:C:257:SER:HB2	2:C:501:NDP:P2B	2.43	0.58
1:C:17:ASP:O	1:C:21:LEU:HD23	2.03	0.58
1:B:157:ASN:HB3	3:C:673:HOH:O	2.03	0.58
1:C:393:ALA:CB	3:C:620:HOH:O	2.43	0.58
1:A:70:LEU:HB3	3:A:607:HOH:O	2.04	0.58
1:D:239[B]:SER:HB2	1:D:387:THR:HG21	1.86	0.58
1:D:49:ASP:O	1:D:52:LYS:HG2	2.04	0.58
1:B:224:GLN:O	3:B:602:HOH:O	2.17	0.57
1:B:271:GLU:OE2	3:B:622:HOH:O	2.17	0.57
1:D:276:ARG:NE	2:D:501:NDP:O3X	2.34	0.57
1:A:286:ASP:OD1	1:A:288:HIS:N	2.36	0.57
1:B:45:SER:O	1:B:405:HIS:NE2	2.30	0.57
1:C:301:LYS:CE	3:C:612:HOH:O	2.43	0.57
1:A:338:THR:HG21	1:C:351:ASN:ND2	2.19	0.57
1:A:436:LEU:HD23	3:A:620:HOH:O	2.04	0.57
1:B:93:THR:HG22	3:B:712:HOH:O	2.05	0.57
1:D:429:PRO:HA	1:D:432:ARG:HB3	1.86	0.57
1:B:133:ASN:CB	3:B:707:HOH:O	2.45	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:43:LEU:O	3:D:626:HOH:O	2.16	0.57
1:B:28:ARG:NE	1:B:56:LEU:HD21	2.20	0.57
1:D:126:LEU:HD11	1:D:180:LEU:HD22	1.87	0.57
1:A:282:ARG:O	1:A:294:GLU:HG3	2.05	0.57
1:A:42:GLY:C	1:A:43:LEU:HD23	2.29	0.56
1:B:32:LEU:HD21	1:B:86:SER:HA	1.86	0.56
1:C:30:VAL:O	1:C:116:VAL:HA	2.04	0.56
1:A:161:MET:HE1	1:C:333:GLY:CA	2.36	0.56
1:A:235:GLY:O	1:A:239[B]:SER:OG	2.20	0.56
1:A:350:ASP:HB3	1:A:353:TYR:CD1	2.40	0.56
1:B:96:PHE:HZ	1:B:166:ARG:NE	2.03	0.56
1:D:234:SER:HB3	1:D:261:LYS:HB3	1.87	0.56
1:D:142:VAL:CG2	3:D:613:HOH:O	2.44	0.56
1:D:350:ASP:OD2	1:D:352:ARG:NH2	2.38	0.56
1:C:282:ARG:O	1:C:294:GLU:HG3	2.06	0.56
1:C:392:TYR:CD1	1:C:400:ALA:HB2	2.41	0.56
1:B:314:VAL:HG21	1:B:326:SER:HB3	1.86	0.56
1:D:403:ASP:HB2	3:D:603:HOH:O	2.05	0.56
1:C:144:HIS:NE2	1:C:187:GLN:HA	2.20	0.56
1:A:23:LEU:HD13	1:A:61:LEU:HD21	1.87	0.56
1:B:268:MET:O	1:B:438:LYS:HE2	2.06	0.56
1:C:74:MET:HE3	1:C:419:ALA:HB2	1.88	0.56
1:A:80:TYR:HB2	3:A:625:HOH:O	2.05	0.56
1:A:437:THR:CG2	3:A:604:HOH:O	2.54	0.56
1:D:100:TYR:CD1	1:D:103:LEU:HD13	2.41	0.56
1:C:231:TRP:CZ2	1:C:322:THR:HG23	2.41	0.56
1:D:96:PHE:HB2	3:D:679:HOH:O	2.05	0.56
1:A:39:MET:HE2	1:A:40:PHE:CE2	2.41	0.56
1:A:231:TRP:CE3	1:A:255:VAL:HG21	2.41	0.56
1:A:440:ASN:OD1	1:A:440:ASN:N	2.35	0.55
1:B:7:ALA:HB2	1:B:21:LEU:HD11	1.88	0.55
1:C:81:ASN:HB2	1:C:170:PHE:CD2	2.41	0.55
1:B:112:LEU:HD13	1:B:114:TYR:CZ	2.42	0.55
1:A:296:LYS:HZ3	1:A:296:LYS:HB2	1.72	0.55
1:C:255:VAL:CG2	3:C:610:HOH:O	2.54	0.55
1:A:1:MET:HE2	1:A:418:LEU:HD21	1.87	0.55
1:A:141:VAL:HA	3:A:605:HOH:O	2.06	0.55
1:B:231:TRP:HA	1:B:255:VAL:CG1	2.37	0.55
1:A:433:SER:CA	3:A:620:HOH:O	2.52	0.55
1:A:332:LYS:NZ	1:C:165:GLN:HG3	2.22	0.55
1:B:158:ASP:OD2	1:B:368:ASN:ND2	2.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:LEU:O	1:D:127:ARG:NH2	2.40	0.55
1:B:46:ARG:HD2	1:B:406:HIS:CD2	2.41	0.55
1:C:223:LYS:NZ	3:C:635:HOH:O	2.31	0.55
1:A:49:ASP:OD2	1:A:52:LYS:HG2	2.07	0.55
1:C:317:HIS:H	1:C:318:PRO:HD2	1.71	0.55
1:B:369:TYR:CD2	1:C:153:PRO:HA	2.41	0.54
1:D:8:ILE:HD13	1:D:74:MET:HE3	1.89	0.54
1:C:214:LEU:HG	1:C:222:LEU:CD2	2.37	0.54
1:A:51:ARG:HH12	1:A:402:LEU:HD13	1.70	0.54
1:A:331:ARG:HG2	1:A:332:LYS:N	2.21	0.54
1:B:199:GLU:O	1:B:203:PRO:HD3	2.08	0.54
1:C:-1:ARG:N	3:C:641:HOH:O	2.40	0.54
1:B:98:GLU:HG3	1:B:99:ARG:HG2	1.90	0.54
1:D:367:ALA:HB1	1:D:371:GLU:HB2	1.89	0.54
1:D:214:LEU:HD21	1:D:222:LEU:CD1	2.37	0.54
1:A:231:TRP:HA	1:A:255:VAL:HG22	1.90	0.54
1:A:161:MET:HE1	1:C:333:GLY:HA3	1.89	0.54
1:A:234:SER:OG	1:A:261:LYS:HG2	2.07	0.54
1:A:433:SER:HA	3:A:620:HOH:O	2.06	0.54
1:A:146:LEU:HD11	1:A:165:GLN:HB2	1.89	0.54
1:D:45:SER:O	3:D:628:HOH:O	2.19	0.54
1:D:127:ARG:HG3	1:D:127:ARG:HH11	1.73	0.54
1:C:101:GLY:C	1:C:103:LEU:H	2.15	0.54
1:A:-1:ARG:CD	3:A:643:HOH:O	2.54	0.53
1:B:191:LYS:NZ	3:B:611:HOH:O	1.99	0.53
1:D:237:LEU:HD21	1:D:339:CYS:SG	2.48	0.53
1:A:352:ARG:HD3	1:C:345:TYR:CD2	2.43	0.53
1:C:403:ASP:O	1:C:408:LYS:HB2	2.09	0.53
1:A:84:TRP:O	1:A:87:ILE:HG22	2.09	0.53
1:A:317:HIS:H	1:A:318:PRO:CD	2.21	0.53
1:A:418:LEU:HA	3:A:613:HOH:O	1.87	0.53
1:D:72:ALA:HB2	1:D:126:LEU:HD21	1.90	0.53
1:D:167:ILE:HG21	1:D:170:PHE:HD2	1.74	0.53
1:D:432:ARG:NH1	1:D:439:ILE:HG21	2.23	0.53
1:A:203:PRO:O	1:A:206:VAL:HG12	2.07	0.53
1:A:302:ILE:O	1:A:306:THR:HG22	2.09	0.53
1:B:342:THR:HG21	3:B:635:HOH:O	2.08	0.53
1:D:296:LYS:CE	1:D:303:ARG:HH22	2.21	0.53
1:A:19:ALA:C	1:A:127:ARG:HH21	2.16	0.53
1:A:168:TRP:HA	1:A:172:THR:CG2	2.39	0.53
1:A:212:ARG:HG3	1:A:371:GLU:OE2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:GLU:CG	3:B:627:HOH:O	2.37	0.53
1:B:145:CYS:SG	1:B:205:LEU:HD22	2.48	0.53
1:A:36:GLU:HB3	1:A:39:MET:SD	2.48	0.53
1:C:214:LEU:HG	1:C:222:LEU:HD22	1.89	0.53
1:B:48:LYS:NZ	1:B:89:GLU:OE2	2.37	0.53
1:A:40:PHE:HB3	1:A:48:LYS:HD3	1.91	0.53
1:A:310:ASP:HB2	1:A:331:ARG:HB2	1.90	0.53
1:B:225:GLY:C	3:B:602:HOH:O	2.36	0.53
1:C:46:ARG:HE	1:C:406:HIS:CD2	2.27	0.53
1:C:199:GLU:O	1:C:202:SER:HB3	2.09	0.53
1:D:69:ALA:HB2	1:D:185:THR:HG22	1.91	0.52
1:A:64:LEU:C	3:A:602:HOH:O	2.38	0.52
1:C:202:SER:OG	1:C:412:LYS:NZ	2.37	0.52
1:A:437:THR:HG23	3:A:604:HOH:O	2.09	0.52
1:D:276:ARG:NH2	3:D:616:HOH:O	2.32	0.52
1:C:67:GLY:O	1:C:185:THR:HG22	2.09	0.52
1:A:222:LEU:HD11	1:A:245:ALA:HA	1.92	0.52
1:A:411:GLY:HA2	3:A:654:HOH:O	2.10	0.52
1:A:191:LYS:N	3:A:605:HOH:O	2.41	0.52
1:D:340:ALA:CB	2:D:501:NDP:H2D	2.40	0.52
1:C:316:GLU:HG3	1:C:317:HIS:N	2.23	0.52
1:B:317:HIS:CG	1:B:318:PRO:HD3	2.44	0.52
1:B:420:PRO:HG2	1:B:424:LEU:HD11	1.91	0.52
1:A:365:HIS:HB3	1:C:358:LEU:HD11	1.92	0.52
1:C:72:ALA:HB3	3:C:615:HOH:O	2.09	0.52
1:A:48:LYS:HD2	1:A:87:ILE:HG13	1.92	0.52
1:A:159:THR:HG21	1:D:156:HIS:ND1	2.25	0.52
1:D:80:TYR:HE1	2:D:501:NDP:O2D	1.93	0.52
1:C:97:LEU:HD11	1:C:115:HIS:CG	2.45	0.52
1:C:125:VAL:HG12	1:C:137:PRO:HA	1.90	0.52
1:A:77:SER:O	1:A:177:LEU:HD21	2.10	0.52
1:D:261:LYS:HA	1:D:264:ILE:HG22	1.91	0.51
1:C:192:PRO:HB2	1:C:195:LEU:HD23	1.93	0.51
1:B:231:TRP:HE3	1:B:255:VAL:HG11	1.71	0.51
1:C:443:ARG:O	1:C:443:ARG:HG2	2.09	0.51
1:B:62:PRO:HG2	1:B:182:LEU:HD13	1.92	0.51
1:D:282:ARG:O	1:D:294:GLU:HG3	2.10	0.51
1:A:327:VAL:HG22	1:A:336:ILE:HD13	1.93	0.51
1:A:38:ARG:CZ	1:A:41:ASP:OD1	2.59	0.51
1:B:28:ARG:CZ	1:B:56:LEU:CD2	2.88	0.51
1:D:168:TRP:HA	1:D:172:THR:OG1	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:CYS:O	1:B:168:TRP:N	2.33	0.51
1:C:189:LEU:O	3:C:606:HOH:O	2.19	0.51
1:B:93:THR:HB	3:B:710:HOH:O	2.11	0.51
1:B:119:SER:OG	1:B:168:TRP:O	2.21	0.51
1:D:321:GLU:OE2	2:D:501:NDP:N6A	2.43	0.51
1:A:38:ARG:NH2	1:A:41:ASP:OD1	2.44	0.51
1:A:144:HIS:O	1:A:187:GLN:HB3	2.11	0.51
1:A:399:GLN:NE2	1:A:402:LEU:HD22	2.26	0.51
1:C:144:HIS:HB2	1:C:366:PHE:CZ	2.46	0.51
1:C:258:SER:CB	1:C:261:LYS:HD3	2.38	0.51
1:D:66:PRO:HB3	1:D:130:ALA:HB3	1.93	0.51
1:C:350:ASP:OD2	3:C:628:HOH:O	2.19	0.51
1:D:296:LYS:NZ	1:D:303:ARG:HH22	2.09	0.51
1:C:286:ASP:OD2	1:C:289:HIS:HB2	2.11	0.51
1:D:82:THR:HA	1:D:85:SER:HB2	1.93	0.50
1:D:330:THR:O	1:D:359:LYS:NZ	2.31	0.50
1:A:370:ARG:HD2	1:D:157:ASN:HB2	1.92	0.50
1:D:104:SER:OG	1:D:107:THR:HG23	2.11	0.50
1:C:118:GLY:HA3	1:C:177:LEU:H	1.76	0.50
1:C:12:ASP:OD2	3:C:627:HOH:O	2.19	0.50
1:C:78:VAL:N	3:C:604:HOH:O	2.39	0.50
1:A:192:PRO:HB3	1:A:380:ALA:HA	1.92	0.50
1:B:323:PHE:HZ	1:B:361:ILE:HD13	1.75	0.50
1:D:145:CYS:HB2	1:D:167:ILE:HD11	1.93	0.50
1:C:95:GLY:HA3	3:C:736:HOH:O	2.12	0.50
1:C:190:PRO:C	3:C:606:HOH:O	2.55	0.50
1:C:205:LEU:HD21	2:C:501:NDP:H5N	1.94	0.50
1:C:258:SER:OG	1:C:260:ARG:NH1	2.44	0.50
1:D:68:GLU:O	1:D:129:GLY:N	2.40	0.50
1:D:132:VAL:HG11	1:D:185:THR:HG21	1.94	0.50
1:B:202:SER:HB3	1:B:388:LEU:HD13	1.93	0.50
1:B:311:VAL:HG22	1:B:313:ILE:H	1.77	0.50
1:C:259:PRO:HG2	1:C:260:ARG:HD3	1.94	0.50
1:A:179:GLN:O	1:A:180:LEU:HD23	2.12	0.49
1:D:97:LEU:O	1:D:111:ASP:HB2	2.12	0.49
1:D:403:ASP:CB	3:D:603:HOH:O	2.59	0.49
1:C:36:GLU:CD	1:C:55:HIS:HE2	2.20	0.49
1:C:100:TYR:HD2	1:C:166:ARG:NH1	2.10	0.49
1:B:256:VAL:O	1:B:276:ARG:N	2.39	0.49
1:C:351:ASN:HB3	1:C:355:TRP:HZ3	1.76	0.49
1:A:282:ARG:CZ	3:A:616:HOH:O	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:34:LYS:O	1:D:37:GLU:HG2	2.12	0.49
1:D:237:LEU:N	2:D:501:NDP:O2N	2.45	0.49
1:C:214:LEU:HD13	1:C:315:PHE:CD2	2.45	0.49
1:A:16:ALA:HB3	3:A:618:HOH:O	2.11	0.49
1:D:410:GLN:HB2	3:D:622:HOH:O	2.13	0.49
1:A:365:HIS:NE2	1:C:356:MET:HG2	2.28	0.49
1:D:61:LEU:HD21	1:D:180:LEU:HD12	1.94	0.49
1:C:147:SER:OG	1:C:166:ARG:HB2	2.13	0.49
1:C:193:LYS:HA	3:C:724:HOH:O	2.13	0.49
1:B:202:SER:O	1:B:412:LYS:HD3	2.12	0.49
1:A:80:TYR:CA	3:A:625:HOH:O	2.61	0.49
1:B:96:PHE:CZ	1:B:166:ARG:NE	2.79	0.49
1:C:316:GLU:OE1	1:C:322:THR:HG22	2.11	0.49
1:A:399:GLN:HE22	1:A:402:LEU:HD22	1.77	0.49
1:B:213:GLN:OE1	1:B:339:CYS:HB2	2.11	0.49
1:B:393:ALA:O	3:B:607:HOH:O	2.12	0.49
1:A:377:ARG:NE	1:B:309:GLU:OE2	2.44	0.49
1:B:321:GLU:CD	2:B:501:NDP:H62A	2.21	0.49
1:B:333:GLY:HA2	1:B:358:LEU:O	2.13	0.49
1:C:77:SER:HB2	3:C:604:HOH:O	2.12	0.48
1:A:25:GLU:HA	1:A:61:LEU:HD23	1.95	0.48
1:D:167:ILE:CG2	1:D:170:PHE:HD2	2.26	0.48
1:A:158:ASP:HB3	1:A:161:MET:HG3	1.96	0.48
1:C:246:LEU:CD1	1:C:252:PRO:HG3	2.42	0.48
1:B:159:THR:HG23	1:C:156:HIS:O	2.13	0.48
1:B:257:SER:OG	3:B:604:HOH:O	1.82	0.48
1:B:257:SER:OG	2:B:501:NDP:O1X	2.31	0.48
1:D:74:MET:HG3	1:D:197:TRP:CE2	2.48	0.48
1:D:317:HIS:H	1:D:318:PRO:CD	2.26	0.48
1:C:23:LEU:HD11	1:C:127:ARG:NE	2.29	0.48
1:A:84:TRP:CA	1:A:87:ILE:HG22	2.41	0.48
1:B:291:ASP:OD2	1:B:293:ARG:HG2	2.12	0.48
1:D:73:VAL:O	1:D:418:LEU:HD13	2.13	0.48
1:D:258:SER:CB	1:D:261:LYS:NZ	2.76	0.48
1:A:190:PRO:C	3:A:605:HOH:O	2.42	0.48
1:A:376:ASN:HB3	3:A:652:HOH:O	2.12	0.48
1:B:430:GLU:O	1:B:434:LYS:HG3	2.13	0.48
1:C:298:LEU:HD23	1:C:325:ALA:HB1	1.95	0.48
1:A:283:PHE:O	1:A:291:ASP:N	2.32	0.48
1:A:366:PHE:CE2	2:A:501:NDP:H41N	2.48	0.48
1:D:327:VAL:HG12	1:D:336:ILE:HD13	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:ILE:O	1:A:361:ILE:HA	2.14	0.48
1:B:167:ILE:HG12	1:B:171:GLU:OE2	2.13	0.48
1:B:298:LEU:HD12	3:B:603:HOH:O	2.14	0.48
1:C:149:GLU:OE2	3:C:629:HOH:O	2.20	0.48
1:B:350:ASP:OD2	1:B:352:ARG:HB2	2.14	0.48
1:D:35:GLY:N	3:D:620:HOH:O	2.07	0.48
1:D:80:TYR:HE1	2:D:501:NDP:HO2N	1.56	0.48
1:C:12:ASP:CG	3:C:627:HOH:O	2.56	0.48
1:A:87:ILE:HG23	1:A:89:GLU:HB2	1.95	0.47
1:A:355:TRP:O	1:C:365:HIS:HB2	2.14	0.47
1:A:418:LEU:CD2	3:A:613:HOH:O	2.49	0.47
1:B:166:ARG:HG3	1:B:166:ARG:HH11	1.79	0.47
1:B:191:LYS:CE	3:B:611:HOH:O	2.55	0.47
1:C:341:SER:OG	1:C:347:HIS:NE2	2.38	0.47
1:D:266:ARG:HG2	1:D:273:ILE:HD12	1.95	0.47
1:D:332:LYS:NZ	1:C:370:ARG:HH12	2.11	0.47
1:A:-1:ARG:NE	3:A:643:HOH:O	2.47	0.47
1:B:189:LEU:HD13	1:B:376:ASN:HB2	1.96	0.47
1:C:78:VAL:HB	1:C:413:VAL:HG23	1.95	0.47
1:C:167:ILE:N	1:C:171:GLU:OE1	2.43	0.47
2:A:501:NDP:O2D	3:A:629:HOH:O	2.19	0.47
1:B:217:ARG:HG3	3:B:700:HOH:O	2.13	0.47
1:C:36:GLU:HB3	1:C:39:MET:HE2	1.95	0.47
1:B:61:LEU:HD12	1:B:62:PRO:O	2.14	0.47
1:C:350:ASP:HB3	1:C:353:TYR:HD2	1.79	0.47
1:A:38:ARG:C	1:A:40:PHE:N	2.65	0.47
1:A:347:HIS:O	1:C:348:GLN:HA	2.15	0.47
1:D:120:ASP:OD1	3:D:629:HOH:O	2.20	0.47
1:C:102:ARG:HA	1:C:108:ALA:CB	2.44	0.47
1:C:118:GLY:HA3	1:C:177:LEU:N	2.29	0.47
1:C:191:LYS:O	1:C:193:LYS:HE3	2.14	0.47
1:B:2:GLN:N	3:B:619:HOH:O	2.07	0.47
1:B:231:TRP:O	1:B:317:HIS:HE1	1.97	0.47
1:C:255:VAL:N	3:C:610:HOH:O	2.41	0.47
1:A:280:GLY:O	1:A:297:ARG:NH2	2.47	0.47
1:D:264:ILE:HD11	1:D:442:PHE:CG	2.50	0.47
1:C:59:VAL:CG2	3:C:618:HOH:O	2.63	0.47
1:C:205:LEU:CD2	2:C:501:NDP:H5N	2.44	0.47
1:A:46:ARG:NE	3:A:645:HOH:O	2.48	0.47
1:A:80:TYR:CZ	2:A:501:NDP:O2A	2.68	0.47
1:A:338:THR:CG2	1:C:351:ASN:HD21	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:SER:CB	1:C:355:TRP:CH2	2.89	0.47
1:B:154:ASP:O	1:C:370:ARG:NH1	2.39	0.47
1:B:352:ARG:HB3	1:B:356:MET:HE2	1.96	0.47
1:D:407:ASN:C	1:D:409:HIS:H	2.23	0.47
1:A:33:HIS:NE2	1:A:57:ASP:OD2	2.37	0.47
1:D:280:GLY:N	3:D:630:HOH:O	2.33	0.47
1:C:377:ARG:HD3	3:C:728:HOH:O	2.15	0.47
1:A:350:ASP:OD2	1:A:352:ARG:NH1	2.47	0.46
1:A:110:HIS:ND1	1:A:172:THR:OG1	2.49	0.46
1:A:204:GLY:N	3:A:638:HOH:O	2.42	0.46
1:B:348:GLN:HA	1:D:347:HIS:O	2.14	0.46
1:C:121:LEU:HG	1:C:143:ALA:HB3	1.97	0.46
1:C:320:ARG:CZ	3:C:613:HOH:O	2.53	0.46
1:B:31:THR:HA	1:B:115:HIS:O	2.15	0.46
1:B:228:VAL:HG13	1:B:313:ILE:HB	1.97	0.46
1:D:291:ASP:OD2	1:D:293:ARG:HB2	2.15	0.46
1:B:83:VAL:O	1:B:87:ILE:HG12	2.16	0.46
1:C:135:TRP:HA	3:C:640:HOH:O	2.16	0.46
1:C:144:HIS:CE1	1:C:187:GLN:HA	2.51	0.46
1:A:51:ARG:HG3	1:A:398:GLY:CA	2.45	0.46
1:A:306:THR:HG21	1:A:309:GLU:O	2.14	0.46
1:B:19:ALA:O	1:B:127:ARG:NH2	2.49	0.46
1:B:226:ASP:OD1	3:B:602:HOH:O	2.20	0.46
1:D:36:GLU:OE1	1:D:55:HIS:HE1	1.97	0.46
1:C:-1:ARG:NH2	3:C:649:HOH:O	2.47	0.46
1:C:193:LYS:HD3	3:C:724:HOH:O	2.16	0.46
1:A:369:TYR:HB3	1:D:156:HIS:CG	2.51	0.46
1:D:340:ALA:HB3	2:D:501:NDP:C2D	2.45	0.46
1:D:145:CYS:O	1:D:167:ILE:HD12	2.16	0.46
1:D:300:GLY:CA	1:D:303:ARG:HH21	2.27	0.46
1:C:317:HIS:N	1:C:318:PRO:CD	2.78	0.46
1:C:353:TYR:O	1:C:357:SER:HB2	2.16	0.46
1:A:48:LYS:NZ	1:A:89:GLU:OE1	2.48	0.46
1:A:352:ARG:O	1:A:356:MET:HB2	2.16	0.46
1:B:151:GLU:OE1	1:C:134:ALA:N	2.37	0.46
1:B:357:SER:HB2	1:B:359:LYS:HE3	1.98	0.46
1:C:116:VAL:HG13	1:C:175:GLY:H	1.80	0.46
1:C:296:LYS:HG2	1:C:328:TYR:CZ	2.51	0.46
1:C:88:PHE:HD1	3:C:671:HOH:O	1.99	0.46
1:A:303:ARG:HA	1:A:306:THR:CG2	2.46	0.46
1:B:96:PHE:CE2	1:B:171:GLU:HB3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:VAL:CG2	1:B:237:LEU:HD12	2.45	0.46
1:C:206:VAL:HG21	1:C:236:GLY:C	2.40	0.46
1:A:83:VAL:HG21	1:A:404:VAL:HB	1.98	0.45
1:A:397:THR:O	1:A:400:ALA:HB3	2.16	0.45
1:B:196:THR:HG23	1:B:199:GLU:H	1.81	0.45
1:B:345:TYR:CE1	1:B:346:MET:HE2	2.51	0.45
1:A:275:ASP:HB3	1:A:278:ALA:HB3	1.99	0.45
1:B:191:LYS:NZ	1:B:195:LEU:O	2.48	0.45
1:C:406:HIS:C	1:C:408:LYS:H	2.24	0.45
1:A:74:MET:HG3	1:A:197:TRP:CD2	2.51	0.45
2:A:501:NDP:H2N	3:A:629:HOH:O	2.15	0.45
1:D:43:LEU:HB3	3:D:601:HOH:O	1.92	0.45
1:D:237:LEU:HD22	2:D:501:NDP:H51N	1.97	0.45
1:C:198:GLU:H	1:C:198:GLU:CD	2.21	0.45
1:C:264:ILE:HG13	1:C:268:MET:HE2	1.98	0.45
1:A:228:VAL:HG23	1:A:228:VAL:O	2.16	0.45
1:B:103:LEU:HD23	1:B:107:THR:HG21	1.98	0.45
1:B:227:ASN:O	1:B:311:VAL:HG23	2.16	0.45
1:D:316:GLU:OE2	1:D:317:HIS:N	2.50	0.45
1:C:100:TYR:HD2	1:C:166:ARG:CZ	2.30	0.45
1:A:32:LEU:HD22	1:A:117:LEU:HD21	1.99	0.45
1:A:68:GLU:OE1	1:A:182:LEU:HD11	2.17	0.45
1:A:378:LEU:HA	1:A:378:LEU:HD23	1.76	0.45
1:B:346:MET:HE1	1:D:350:ASP:HB2	1.98	0.45
1:B:428:ASP:OD1	1:B:431:LEU:HB3	2.17	0.45
1:D:125:VAL:HG23	1:D:137:PRO:HA	1.98	0.45
1:D:212:ARG:HD2	1:D:367:ALA:HA	1.98	0.45
1:D:262:ALA:O	1:D:266:ARG:HG3	2.16	0.45
1:C:292:PRO:N	3:C:617:HOH:O	2.47	0.45
1:A:8:ILE:HG12	1:A:124:VAL:HG21	1.98	0.45
1:D:279:GLU:HB3	3:D:621:HOH:O	2.17	0.45
1:C:390:LYS:HE3	1:C:392:TYR:CE1	2.51	0.45
1:B:99:ARG:HD3	1:B:102:ARG:HB2	1.97	0.45
1:B:198:GLU:CD	1:B:198:GLU:H	2.25	0.45
1:D:54:LEU:HD12	1:D:398:GLY:HA2	1.98	0.45
1:D:308:GLY:C	3:D:643:HOH:O	2.59	0.45
1:C:144:HIS:HB2	1:C:366:PHE:HZ	1.82	0.45
1:A:366:PHE:CD2	2:A:501:NDP:H41N	2.51	0.45
1:B:1:MET:HE3	1:B:179:GLN:OE1	2.17	0.45
1:B:11:GLY:O	1:B:13:ALA:N	2.50	0.45
1:B:355:TRP:CZ3	1:D:340:ALA:HA	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:339:CYS:O	2:D:501:NDP:N7N	2.33	0.45
1:C:74:MET:HE2	1:C:197:TRP:NE1	2.32	0.45
1:A:285:LYS:HE3	1:A:289:HIS:HB3	1.99	0.44
1:B:160:MET:HE2	1:B:365:HIS:O	2.17	0.44
1:B:192:PRO:HB2	1:B:195:LEU:HD12	1.99	0.44
1:B:229:LEU:N	1:B:311:VAL:HG21	2.26	0.44
1:B:84:TRP:HB3	1:B:89:GLU:O	2.18	0.44
1:B:355:TRP:CD2	1:D:340:ALA:HA	2.51	0.44
1:D:145:CYS:HB2	1:D:167:ILE:CD1	2.47	0.44
1:C:97:LEU:HD11	1:C:115:HIS:CD2	2.53	0.44
1:C:242:THR:O	1:C:246:LEU:HD13	2.18	0.44
1:A:243:GLN:OE1	1:A:385:HIS:N	2.44	0.44
1:B:165:GLN:O	1:B:165:GLN:HG3	2.16	0.44
1:A:160:MET:HE1	1:A:165:GLN:HE22	1.81	0.44
1:A:212:ARG:NH1	1:A:371:GLU:OE1	2.49	0.44
1:B:234:SER:OG	1:B:261:LYS:HD2	2.17	0.44
1:A:32:LEU:CD2	1:A:117:LEU:HD21	2.47	0.44
1:A:179:GLN:HG2	3:A:613:HOH:O	2.18	0.44
1:A:390:LYS:HE3	1:A:392:TYR:CZ	2.52	0.44
1:C:157:ASN:O	1:C:158:ASP:HB2	2.18	0.44
1:A:48:LYS:HD2	1:A:87:ILE:CD1	2.48	0.44
1:A:78:VAL:HG21	1:A:400:ALA:HB1	1.99	0.44
1:B:440:ASN:HA	3:B:673:HOH:O	2.16	0.44
1:C:34:LYS:NZ	1:C:37:GLU:OE2	2.45	0.44
1:C:74:MET:HE2	1:C:197:TRP:CD1	2.52	0.44
1:C:268:MET:HE3	1:C:268:MET:HB2	1.85	0.44
1:A:29:ALA:O	1:A:56:LEU:HA	2.17	0.44
1:B:79:ASN:ND2	3:B:640:HOH:O	2.51	0.44
1:B:80:TYR:HD2	1:B:84:TRP:CH2	2.36	0.44
1:A:400:ALA:O	1:A:404:VAL:HG23	2.18	0.44
1:B:122:ALA:HB2	1:B:201:ALA:HB2	2.00	0.44
1:B:317:HIS:H	1:B:318:PRO:CD	2.31	0.44
1:D:157:ASN:ND2	3:D:654:HOH:O	2.51	0.44
1:C:74:MET:HG3	1:C:197:TRP:CE2	2.52	0.44
1:C:78:VAL:HG21	1:C:400:ALA:HB1	2.00	0.44
1:C:369:TYR:HB2	3:C:665:HOH:O	2.17	0.44
1:A:1:MET:HE3	1:A:1:MET:HA	1.99	0.44
1:A:76:SER:HB3	1:A:177:LEU:HD23	1.99	0.44
1:B:37:GLU:HA	1:B:88:PHE:CD1	2.53	0.44
1:B:191:LYS:HE2	3:B:611:HOH:O	2.18	0.44
1:D:94:PHE:HA	1:D:97:LEU:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:ILE:O	1:A:8:ILE:HD12	2.17	0.43
1:D:409:HIS:CE1	2:D:501:NDP:O1A	2.71	0.43
1:B:260:ARG:NE	3:B:624:HOH:O	2.24	0.43
1:D:32:LEU:HD11	1:D:85:SER:C	2.42	0.43
1:D:231:TRP:NE1	1:D:326:SER:OG	2.49	0.43
1:D:336:ILE:O	1:D:361:ILE:HA	2.17	0.43
1:C:191:LYS:N	3:C:606:HOH:O	2.51	0.43
1:B:28:ARG:HE	1:B:56:LEU:CD2	2.31	0.43
1:C:100:TYR:CD2	1:C:166:ARG:CZ	3.01	0.43
1:C:121:LEU:HD12	1:C:121:LEU:C	2.43	0.43
1:C:233:ALA:HA	1:C:238:GLY:HA3	2.00	0.43
1:A:83:VAL:HG22	1:A:401:ALA:HB1	1.99	0.43
1:B:151:GLU:OE2	1:C:133:ASN:HB3	2.19	0.43
1:D:27:TYR:HB2	1:D:179:GLN:HB2	1.99	0.43
1:D:167:ILE:HD13	1:D:366:PHE:CE1	2.53	0.43
1:D:205:LEU:HD21	1:D:366:PHE:HE2	1.83	0.43
1:C:64:LEU:HD11	1:C:70:LEU:HD13	2.00	0.43
1:C:209:THR:OG1	2:C:501:NDP:H41N	2.18	0.43
1:A:60:PRO:HD2	1:A:114:TYR:OH	2.18	0.43
1:A:121:LEU:HD22	1:A:143:ALA:HB2	2.01	0.43
1:C:205:LEU:HA	1:C:366:PHE:HE1	1.83	0.43
1:B:28:ARG:NH1	1:B:58:ASP:OD2	2.51	0.43
1:B:117:LEU:HB3	3:B:716:HOH:O	2.18	0.43
1:B:193:LYS:HB3	1:B:427:THR:OG1	2.19	0.43
1:C:298:LEU:O	1:C:302:ILE:HG13	2.19	0.43
1:A:5:LEU:HD13	1:A:418:LEU:HB3	2.00	0.43
1:A:14:ALA:O	1:A:17:ASP:HB2	2.19	0.43
1:B:227:ASN:C	1:B:311:VAL:HG23	2.43	0.43
1:D:80:TYR:O	1:D:83:VAL:HB	2.19	0.43
1:D:235:GLY:HA2	1:D:410:GLN:O	2.18	0.43
1:A:362:VAL:HG22	1:C:360:ARG:HG2	2.00	0.43
1:B:120:ASP:OD2	1:B:204:GLY:HA3	2.18	0.43
1:C:306:THR:HB	1:C:309:GLU:HB2	2.00	0.43
1:A:80:TYR:CB	3:A:625:HOH:O	2.64	0.43
1:B:106:LEU:HD23	1:B:106:LEU:HA	1.77	0.43
1:B:314:VAL:CG2	1:B:336:ILE:HG23	2.49	0.43
1:D:8:ILE:HD13	1:D:74:MET:CE	2.49	0.43
1:B:32:LEU:HD13	3:B:712:HOH:O	2.19	0.42
1:A:301:LYS:HA	1:A:301:LYS:HD3	1.79	0.42
1:D:308:GLY:CA	3:D:618:HOH:O	2.59	0.42
1:C:336:ILE:N	1:C:336:ILE:HD12	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:GLU:O	1:A:297:ARG:NH1	2.49	0.42
1:B:287:GLU:O	1:B:288:HIS:ND1	2.52	0.42
1:D:54:LEU:HD21	1:D:86:SER:HB3	2.00	0.42
1:D:100:TYR:CE1	1:D:103:LEU:HD22	2.54	0.42
1:D:158:ASP:OD2	1:D:368:ASN:ND2	2.44	0.42
1:D:309:GLU:CD	1:C:377:ARG:HH21	2.22	0.42
1:C:261:LYS:HG2	1:C:410:GLN:OE1	2.19	0.42
1:A:65:GLY:CA	3:A:602:HOH:O	2.60	0.42
1:A:418:LEU:HD23	3:A:613:HOH:O	2.17	0.42
1:B:54:LEU:HD23	1:B:54:LEU:HA	1.83	0.42
1:D:31:THR:HG23	1:D:114:TYR:HB2	2.01	0.42
1:C:56:LEU:HD11	1:C:394:LEU:HD22	2.00	0.42
1:C:252:PRO:O	1:C:272:ALA:HB3	2.20	0.42
1:A:48:LYS:HB2	1:A:48:LYS:HE3	1.83	0.42
1:A:377:ARG:NH1	3:A:641:HOH:O	2.45	0.42
1:B:95:GLY:O	1:B:98:GLU:HG2	2.19	0.42
1:B:179:GLN:O	1:B:180:LEU:HD23	2.19	0.42
1:D:14:ALA:HB3	1:D:17:ASP:OD2	2.19	0.42
1:C:5:LEU:O	1:C:9:LEU:HG	2.19	0.42
1:C:301:LYS:HD2	1:C:304:GLU:OE1	2.19	0.42
1:C:349:TYR:CE2	1:C:351:ASN:HA	2.54	0.42
1:C:426:VAL:HB	3:C:616:HOH:O	2.20	0.42
1:C:56:LEU:HD13	1:C:56:LEU:HA	1.81	0.42
1:C:323:PHE:O	1:C:327:VAL:HG23	2.20	0.42
1:C:353:TYR:O	1:C:359:LYS:HE3	2.20	0.42
1:A:44:ALA:HB1	1:A:45:SER:H	1.61	0.42
1:C:191:LYS:HA	3:C:606:HOH:O	2.20	0.42
1:B:217:ARG:NH1	3:B:632:HOH:O	2.45	0.42
1:D:35:GLY:C	1:D:37:GLU:H	2.28	0.42
1:D:127:ARG:HG3	1:D:127:ARG:NH1	2.33	0.42
1:D:223:LYS:HD3	3:D:658:HOH:O	2.19	0.42
1:C:321:GLU:O	3:C:630:HOH:O	2.21	0.42
1:A:34:LYS:HE2	1:A:94:PHE:CG	2.55	0.42
1:A:160:MET:HB3	1:C:358:LEU:HD21	2.02	0.42
1:D:56:LEU:HD12	1:D:56:LEU:HA	1.87	0.42
1:C:214:LEU:C	1:C:222:LEU:HD23	2.45	0.42
1:C:303:ARG:HD3	1:C:308:GLY:HA2	2.02	0.42
1:B:333:GLY:CA	1:D:161:MET:HE1	2.36	0.42
1:A:261:LYS:HD3	2:A:501:NDP:O3X	2.20	0.41
1:B:16:ALA:N	3:B:634:HOH:O	2.47	0.41
1:B:220:ALA:HB1	1:B:313:ILE:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:330:THR:CG2	1:B:359:LYS:HD3	2.49	0.41
1:D:46:ARG:HG3	1:D:405:HIS:CE1	2.55	0.41
1:C:74:MET:HG3	1:C:197:TRP:CD2	2.55	0.41
1:C:161:MET:HE2	1:C:161:MET:HB2	1.78	0.41
1:A:284:TRP:HA	1:A:290:GLN:HA	2.02	0.41
1:D:56:LEU:CD2	1:D:395:GLU:HG3	2.49	0.41
1:B:157:ASN:HB2	1:C:370:ARG:HD3	2.02	0.41
1:D:74:MET:HG3	1:D:197:TRP:CD2	2.56	0.41
1:D:368:ASN:OD1	1:D:371:GLU:HG3	2.20	0.41
1:A:256:VAL:O	1:A:276:ARG:N	2.52	0.41
1:B:352:ARG:HD3	1:D:345:TYR:CG	2.56	0.41
1:D:29:ALA:HA	1:D:178:ALA:HA	2.02	0.41
1:D:161:MET:HE3	1:D:161:MET:HB3	1.36	0.41
1:D:203:PRO:O	1:D:204:GLY:C	2.63	0.41
1:A:116:VAL:N	1:A:173:ASN:OD1	2.52	0.41
1:A:345:TYR:CD2	1:C:352:ARG:HD3	2.55	0.41
1:B:206:VAL:HG22	1:B:237:LEU:HD12	2.02	0.41
1:C:73:VAL:HB	1:C:178:ALA:O	2.20	0.41
1:A:47:ASP:CG	3:A:630:HOH:O	2.63	0.41
1:A:406:HIS:O	1:A:407:ASN:C	2.62	0.41
1:B:154:ASP:HB3	1:D:332:LYS:NZ	2.34	0.41
1:B:177:LEU:HD23	1:B:177:LEU:HA	1.92	0.41
1:D:66:PRO:HA	1:D:130:ALA:H	1.84	0.41
1:C:40:PHE:CE1	1:C:53[B]:SER:HB3	2.56	0.41
1:C:291:ASP:CA	3:C:617:HOH:O	2.55	0.41
1:C:314:VAL:CG1	1:C:336:ILE:HG13	2.51	0.41
1:A:14:ALA:N	1:A:17:ASP:OD2	2.42	0.41
1:A:153:PRO:HG3	1:D:369:TYR:CZ	2.56	0.41
1:A:156:HIS:ND1	1:D:159:THR:HG21	2.35	0.41
1:B:228:VAL:HA	1:B:313:ILE:O	2.20	0.41
1:D:145:CYS:SG	1:D:167:ILE:HD11	2.60	0.41
1:D:308:GLY:HA2	3:D:618:HOH:O	2.19	0.41
1:C:320:ARG:NH2	3:C:613:HOH:O	2.54	0.41
1:A:209:THR:OG1	2:A:501:NDP:O7N	2.39	0.41
1:D:214:LEU:HD11	1:D:228:VAL:HG11	2.03	0.41
1:A:73:VAL:HB	1:A:178:ALA:O	2.21	0.41
1:A:82:THR:N	3:A:623:HOH:O	2.36	0.41
1:A:323:PHE:CZ	1:A:361:ILE:HD13	2.56	0.41
1:B:99:ARG:HA	1:B:102:ARG:HB2	2.03	0.41
1:B:412:LYS:HE2	1:B:412:LYS:HB3	1.62	0.41
1:D:52:LYS:HE2	1:D:52:LYS:HB3	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:99:ARG:CA	3:D:608:HOH:O	2.47	0.41
1:C:59:VAL:HG11	1:C:114:TYR:CE1	2.56	0.41
1:C:296:LYS:HG2	1:C:328:TYR:OH	2.21	0.41
1:A:135:TRP:CD1	1:A:185:THR:HB	2.56	0.41
1:A:349:TYR:HE2	1:A:351:ASN:HA	1.86	0.41
1:B:48:LYS:HB2	1:B:87:ILE:HD12	2.02	0.41
1:B:421:ARG:HG3	1:B:422:GLU:O	2.21	0.41
1:B:431:LEU:HD12	1:B:435:HIS:CD2	2.56	0.41
1:A:279:GLU:O	1:A:297:ARG:NH2	2.51	0.40
1:A:373:PHE:CE2	3:D:725:HOH:O	2.64	0.40
1:A:401:ALA:O	1:A:405:HIS:HB2	2.22	0.40
1:B:99:ARG:HD3	1:B:99:ARG:HA	1.74	0.40
1:D:28:ARG:HG2	1:D:56:LEU:HD11	2.03	0.40
1:A:28:ARG:HA	1:A:57:ASP:O	2.21	0.40
1:A:228:VAL:HA	1:A:313:ILE:O	2.21	0.40
1:A:286:ASP:OD1	1:A:289:HIS:N	2.51	0.40
1:B:30:VAL:HA	1:B:55:HIS:O	2.21	0.40
1:D:442:PHE:HB2	3:D:657:HOH:O	2.20	0.40
1:C:327:VAL:O	1:C:359:LYS:NZ	2.50	0.40
1:C:395:GLU:OE2	1:C:395:GLU:N	2.54	0.40
1:A:40:PHE:HB3	1:A:48:LYS:CD	2.51	0.40
1:A:168:TRP:NE1	1:A:174:PHE:O	2.52	0.40
1:A:227:ASN:ND2	1:A:309:GLU:OE1	2.48	0.40
1:A:357:SER:CB	1:A:359:LYS:HE3	2.48	0.40
1:B:195:LEU:HD22	1:B:386:PRO:HG3	2.03	0.40
1:B:352:ARG:NH1	1:D:345:TYR:CE1	2.90	0.40
1:D:264:ILE:HD11	1:D:442:PHE:HA	2.04	0.40
1:C:25:GLU:O	1:C:60:PRO:HA	2.21	0.40
1:C:377:ARG:O	1:C:381:LYS:HG2	2.21	0.40
1:A:34:LYS:HA	1:A:94:PHE:CE2	2.56	0.40
1:A:255:VAL:HA	1:A:274:ILE:O	2.22	0.40
1:A:389:SER:OG	1:A:442:PHE:HB3	2.21	0.40
1:B:67:GLY:O	1:B:185:THR:HG22	2.21	0.40
1:C:318:PRO:HB2	1:C:322:THR:HB	2.03	0.40
1:A:48:LYS:HD2	1:A:87:ILE:CG1	2.51	0.40
1:A:48:LYS:HD2	1:A:87:ILE:HD11	2.03	0.40
1:A:434:LYS:HE3	1:A:434:LYS:HB2	1.96	0.40
1:A:436:LEU:CB	3:A:620:HOH:O	2.60	0.40
1:B:167:ILE:HD11	1:B:170:PHE:CD2	2.56	0.40
1:D:83:VAL:O	1:D:86:SER:OG	2.33	0.40
1:C:189:LEU:HD13	1:C:376:ASN:CB	2.51	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	444/445 (100%)	408 (92%)	32 (7%)	4 (1%)	14	10
1	B	445/445 (100%)	418 (94%)	21 (5%)	6 (1%)	9	6
1	C	445/445 (100%)	424 (95%)	15 (3%)	6 (1%)	9	6
1	D	445/445 (100%)	401 (90%)	37 (8%)	7 (2%)	7	4
All	All	1779/1780 (100%)	1651 (93%)	105 (6%)	23 (1%)	9	6

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	ALA
1	A	317	HIS
1	B	12	ASP
1	B	317	HIS
1	D	103	LEU
1	D	317	HIS
1	C	103	LEU
1	C	158	ASP
1	C	317	HIS
1	C	366	PHE
1	A	158	ASP
1	B	103	LEU
1	D	10	SER
1	D	204	GLY
1	C	101	GLY
1	C	102	ARG
1	D	37	GLU
1	D	408	LYS
1	A	407	ASN
1	B	13	ALA

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Mol	Chain	Res	Type
1	B	90	PRO
1	B	420	PRO
1	D	11	GLY

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	358/357 (100%)	346 (97%)	12 (3%)	32	35
1	B	359/357 (101%)	357 (99%)	2 (1%)	78	86
1	C	359/357 (101%)	358 (100%)	1 (0%)	86	91
1	D	359/357 (101%)	358 (100%)	1 (0%)	86	91
All	All	1435/1428 (100%)	1419 (99%)	16 (1%)	65	74

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASP
1	A	32	LEU
1	A	37	GLU
1	A	41	ASP
1	A	46	ARG
1	A	48	LYS
1	A	49	ASP
1	A	81	ASN
1	A	104	SER
1	A	410	GLN
1	A	436	LEU
1	A	437	THR
1	B	53[A]	SER
1	B	53[B]	SER
1	D	160	MET
1	C	165	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	GLN
1	A	224	GLN
1	B	2	GLN
1	B	79	ASN
1	B	115	HIS
1	B	207	ASN
1	B	317	HIS
1	B	410	GLN
1	D	55	HIS
1	D	351	ASN
1	C	2	GLN
1	C	194	HIS
1	C	348	GLN
1	C	351	ASN
1	C	385	HIS
1	C	399	GLN
1	C	435	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	NDP	B	501	-	51,52,52	4.00	26 (50%)	71,80,80	1.70	15 (21%)
2	NDP	D	501	-	51,52,52	4.05	22 (43%)	71,80,80	2.03	18 (25%)
2	NDP	C	501	-	51,52,52	3.86	22 (43%)	71,80,80	1.98	19 (26%)
2	NDP	A	501	-	51,52,52	4.00	24 (47%)	71,80,80	1.79	16 (22%)

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	NDP	B	501	-	3/3/14/17	9/34/77/77	0/5/5/5
2	NDP	D	501	-	3/3/14/17	12/34/77/77	0/5/5/5
2	NDP	C	501	-	3/3/14/17	15/34/77/77	0/5/5/5
2	NDP	A	501	-	3/3/14/17	9/34/77/77	0/5/5/5

All (94) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NDP	PA-O3	10.29	1.70	1.59
2	D	501	NDP	PA-O3	10.08	1.70	1.59
2	A	501	NDP	O4B-C1B	9.37	1.63	1.42
2	B	501	NDP	PA-O3	9.31	1.69	1.59
2	C	501	NDP	O4D-C1D	9.08	1.63	1.42
2	D	501	NDP	O4B-C1B	8.91	1.62	1.42
2	B	501	NDP	O4B-C4B	-8.90	1.25	1.45
2	C	501	NDP	O4B-C1B	8.88	1.62	1.42
2	B	501	NDP	O4D-C1D	8.83	1.62	1.42
2	A	501	NDP	O4D-C1D	8.77	1.62	1.42
2	D	501	NDP	O4D-C1D	8.62	1.61	1.42
2	B	501	NDP	O4B-C1B	8.43	1.61	1.42
2	D	501	NDP	C2D-C1D	-8.41	1.27	1.53
2	A	501	NDP	C2B-C1B	-8.26	1.32	1.53
2	C	501	NDP	O4B-C4B	-8.25	1.26	1.45
2	C	501	NDP	O4D-C4D	-8.21	1.26	1.45
2	D	501	NDP	C2B-C1B	-8.13	1.33	1.53
2	B	501	NDP	PN-O3	8.11	1.68	1.59
2	B	501	NDP	C2B-C1B	-8.03	1.33	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	NDP	C2B-C1B	-8.01	1.33	1.53
2	C	501	NDP	P2B-O2B	7.93	1.73	1.59
2	D	501	NDP	O4B-C4B	-7.89	1.27	1.45
2	A	501	NDP	O4D-C4D	-7.78	1.27	1.45
2	B	501	NDP	O4D-C4D	-7.75	1.27	1.45
2	A	501	NDP	O4B-C4B	-7.75	1.27	1.45
2	D	501	NDP	PN-O3	7.59	1.67	1.59
2	D	501	NDP	P2B-O2B	7.57	1.72	1.59
2	D	501	NDP	O4D-C4D	-7.52	1.28	1.45
2	C	501	NDP	PA-O3	7.41	1.67	1.59
2	A	501	NDP	PN-O3	7.20	1.67	1.59
2	A	501	NDP	C2D-C1D	-7.19	1.30	1.53
2	B	501	NDP	P2B-O2B	7.16	1.72	1.59
2	A	501	NDP	P2B-O2B	6.94	1.71	1.59
2	B	501	NDP	C2D-C1D	-6.73	1.32	1.53
2	C	501	NDP	C2D-C1D	-6.52	1.33	1.53
2	D	501	NDP	C4N-C3N	-6.23	1.38	1.50
2	C	501	NDP	PN-O3	6.05	1.66	1.59
2	A	501	NDP	C4N-C3N	-5.94	1.38	1.50
2	C	501	NDP	C4N-C3N	-5.67	1.39	1.50
2	B	501	NDP	C4N-C3N	-5.54	1.39	1.50
2	C	501	NDP	C6A-N6A	4.98	1.46	1.34
2	A	501	NDP	C6A-N6A	4.91	1.46	1.34
2	D	501	NDP	C6A-N6A	4.84	1.46	1.34
2	B	501	NDP	C6A-N6A	4.61	1.46	1.34
2	D	501	NDP	C4N-C5N	-3.95	1.38	1.49
2	D	501	NDP	C5A-N7A	-3.86	1.32	1.39
2	A	501	NDP	C4N-C5N	-3.75	1.39	1.49
2	C	501	NDP	C8A-N9A	-3.74	1.31	1.37
2	B	501	NDP	C8A-N9A	-3.59	1.31	1.37
2	D	501	NDP	C7N-N7N	3.55	1.43	1.33
2	B	501	NDP	O2D-C2D	3.50	1.51	1.43
2	B	501	NDP	C7N-N7N	3.49	1.43	1.33
2	C	501	NDP	C7N-C3N	3.48	1.56	1.48
2	A	501	NDP	O2D-C2D	3.46	1.51	1.43
2	A	501	NDP	O3D-C3D	-3.45	1.34	1.43
2	C	501	NDP	C7N-N7N	3.35	1.43	1.33
2	A	501	NDP	C7N-N7N	3.34	1.43	1.33
2	B	501	NDP	C7N-C3N	3.34	1.55	1.48
2	D	501	NDP	O2B-C2B	3.30	1.55	1.44
2	D	501	NDP	C7N-C3N	3.25	1.55	1.48
2	C	501	NDP	C5A-N7A	-3.24	1.33	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NDP	O3D-C3D	-3.24	1.34	1.43
2	B	501	NDP	C4N-C5N	-3.21	1.40	1.49
2	C	501	NDP	C4N-C5N	-3.17	1.40	1.49
2	A	501	NDP	C8A-N9A	-3.14	1.32	1.37
2	D	501	NDP	O3D-C3D	-3.13	1.35	1.43
2	A	501	NDP	PN-O5D	3.13	1.71	1.59
2	C	501	NDP	O3D-C3D	-3.09	1.35	1.43
2	B	501	NDP	O2B-C2B	3.08	1.54	1.44
2	D	501	NDP	C6N-N1N	-3.07	1.29	1.37
2	A	501	NDP	O7N-C7N	-3.07	1.17	1.24
2	D	501	NDP	O2D-C2D	3.04	1.50	1.43
2	B	501	NDP	PN-O5D	2.99	1.71	1.59
2	C	501	NDP	O2D-C2D	2.94	1.50	1.43
2	A	501	NDP	C5A-N7A	-2.87	1.33	1.39
2	B	501	NDP	C5A-N7A	-2.83	1.33	1.39
2	C	501	NDP	PN-O5D	2.80	1.70	1.59
2	D	501	NDP	O7N-C7N	-2.73	1.18	1.24
2	D	501	NDP	C8A-N9A	-2.70	1.33	1.37
2	C	501	NDP	O7N-C7N	-2.69	1.18	1.24
2	C	501	NDP	O2B-C2B	2.66	1.53	1.44
2	B	501	NDP	O7N-C7N	-2.59	1.18	1.24
2	B	501	NDP	C4A-N9A	-2.55	1.32	1.37
2	C	501	NDP	C6N-N1N	-2.42	1.31	1.37
2	B	501	NDP	C1B-N9A	-2.42	1.39	1.46
2	A	501	NDP	C3B-C4B	2.40	1.59	1.53
2	B	501	NDP	C5D-C4D	2.39	1.58	1.51
2	A	501	NDP	C7N-C3N	2.33	1.53	1.48
2	A	501	NDP	C6N-N1N	-2.28	1.31	1.37
2	B	501	NDP	O3B-C3B	-2.24	1.37	1.43
2	A	501	NDP	O2B-C2B	2.23	1.51	1.44
2	B	501	NDP	C6N-N1N	-2.21	1.31	1.37
2	A	501	NDP	C5D-C4D	2.15	1.58	1.51
2	D	501	NDP	PN-O5D	2.03	1.67	1.59

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	NDP	C5A-C4A-N3A	-6.24	118.12	126.72
2	C	501	NDP	C5A-C4A-N3A	-5.80	118.73	126.72
2	A	501	NDP	C5A-C4A-N3A	-5.68	118.89	126.72
2	D	501	NDP	O5D-C5D-C4D	-5.26	91.07	108.99
2	B	501	NDP	C2B-C1B-N9A	-4.75	105.94	113.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NDP	N3A-C2A-N1A	-4.72	121.43	128.58
2	C	501	NDP	N3A-C2A-N1A	-4.60	121.62	128.58
2	A	501	NDP	C2B-C1B-N9A	-4.59	106.20	113.75
2	B	501	NDP	C5A-C4A-N3A	-4.34	120.74	126.72
2	B	501	NDP	N3A-C2A-N1A	-4.22	122.19	128.58
2	D	501	NDP	O4D-C1D-N1N	4.22	116.13	108.08
2	D	501	NDP	N3A-C4A-N9A	4.14	134.21	127.17
2	D	501	NDP	C6N-N1N-C2N	4.12	123.72	119.32
2	C	501	NDP	N3A-C4A-N9A	4.05	134.05	127.17
2	C	501	NDP	C2D-C1D-N1N	4.03	123.22	113.31
2	D	501	NDP	O4D-C1D-C2D	-3.97	98.11	106.62
2	A	501	NDP	N3A-C4A-N9A	3.85	133.72	127.17
2	A	501	NDP	C2A-N3A-C4A	3.84	121.20	111.83
2	C	501	NDP	C4A-C5A-N7A	-3.68	106.37	110.58
2	A	501	NDP	C4A-C5A-N7A	-3.65	106.41	110.58
2	C	501	NDP	C2A-N3A-C4A	3.64	120.72	111.83
2	B	501	NDP	C4A-C5A-N7A	-3.60	106.47	110.58
2	D	501	NDP	N3A-C2A-N1A	-3.59	123.14	128.58
2	C	501	NDP	O3-PN-O1N	-3.56	99.99	110.70
2	D	501	NDP	C3N-C2N-N1N	-3.56	117.98	123.20
2	D	501	NDP	C4D-O4D-C1D	-3.53	101.67	109.47
2	B	501	NDP	N9A-C8A-N7A	-3.53	108.93	113.94
2	D	501	NDP	C2A-N3A-C4A	3.47	120.30	111.83
2	C	501	NDP	C5A-N7A-C8A	3.45	108.86	103.45
2	D	501	NDP	C1D-N1N-C6N	-3.40	113.58	120.77
2	A	501	NDP	C5A-N7A-C8A	3.35	108.71	103.45
2	B	501	NDP	C5A-N7A-C8A	3.28	108.61	103.45
2	C	501	NDP	C5B-C4B-C3B	-3.23	103.60	115.21
2	A	501	NDP	C3N-C2N-N1N	-3.18	118.53	123.20
2	B	501	NDP	C4A-N9A-C8A	3.15	109.05	105.74
2	B	501	NDP	C2A-N3A-C4A	3.09	119.37	111.83
2	B	501	NDP	N3A-C4A-N9A	3.03	132.32	127.17
2	A	501	NDP	C1D-N1N-C2N	-3.00	116.19	121.14
2	A	501	NDP	C6N-N1N-C2N	2.96	122.49	119.32
2	A	501	NDP	C3B-C2B-C1B	2.95	108.46	102.81
2	C	501	NDP	N9A-C8A-N7A	-2.92	109.79	113.94
2	D	501	NDP	C4A-C5A-N7A	-2.86	107.32	110.58
2	A	501	NDP	N9A-C8A-N7A	-2.80	109.96	113.94
2	C	501	NDP	O2A-PA-O3	2.74	114.68	107.27
2	C	501	NDP	C2D-C3D-C4D	-2.64	97.50	102.61
2	C	501	NDP	O2B-C2B-C3B	-2.62	102.28	111.68
2	A	501	NDP	O4D-C1D-C2D	-2.62	101.02	106.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	NDP	C5A-N7A-C8A	2.61	107.56	103.45
2	B	501	NDP	C4B-O4B-C1B	-2.58	103.76	109.47
2	C	501	NDP	O2D-C2D-C1D	2.58	118.99	110.10
2	C	501	NDP	C4D-O4D-C1D	-2.56	103.82	109.47
2	B	501	NDP	C3N-C2N-N1N	-2.48	119.57	123.20
2	A	501	NDP	O4B-C1B-N9A	2.45	112.79	108.09
2	D	501	NDP	O2B-C2B-C1B	2.41	118.54	110.05
2	C	501	NDP	O3D-C3D-C4D	2.41	118.02	111.08
2	C	501	NDP	C3B-C2B-C1B	2.38	107.37	102.81
2	B	501	NDP	C1D-N1N-C2N	-2.38	117.22	121.14
2	B	501	NDP	O4B-C4B-C3B	-2.33	100.53	105.15
2	C	501	NDP	O4D-C4D-C3D	-2.33	100.54	105.15
2	C	501	NDP	C6N-N1N-C2N	2.24	121.72	119.32
2	A	501	NDP	C5B-C4B-C3B	-2.24	107.15	115.21
2	D	501	NDP	C5B-C4B-C3B	-2.20	107.28	115.21
2	B	501	NDP	C6N-N1N-C2N	2.19	121.66	119.32
2	D	501	NDP	C4B-O4B-C1B	-2.17	104.67	109.47
2	B	501	NDP	O5D-C5D-C4D	-2.15	101.69	108.99
2	D	501	NDP	C6A-C5A-C4A	2.13	120.09	117.18
2	D	501	NDP	O4B-C1B-N9A	2.03	111.98	108.09
2	A	501	NDP	C4D-O4D-C1D	-2.01	105.03	109.47

All (12) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	501	NDP	C2D
2	A	501	NDP	C3D
2	A	501	NDP	C4D
2	B	501	NDP	C2D
2	B	501	NDP	C3D
2	B	501	NDP	C4D
2	D	501	NDP	C2D
2	D	501	NDP	C3D
2	D	501	NDP	C4D
2	C	501	NDP	C2D
2	C	501	NDP	C3D
2	C	501	NDP	C4D

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NDP	C4D-C5D-O5D-PN

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Mol	Chain	Res	Type	Atoms
2	B	501	NDP	C5B-O5B-PA-O1A
2	B	501	NDP	C5B-O5B-PA-O2A
2	B	501	NDP	C5B-O5B-PA-O3
2	B	501	NDP	C5D-O5D-PN-O3
2	B	501	NDP	C5D-O5D-PN-O2N
2	B	501	NDP	O4D-C1D-N1N-C2N
2	B	501	NDP	C2N-C3N-C7N-N7N
2	D	501	NDP	C5B-O5B-PA-O2A
2	D	501	NDP	C5B-O5B-PA-O3
2	D	501	NDP	C1B-C2B-O2B-P2B
2	D	501	NDP	PA-O3-PN-O5D
2	D	501	NDP	C5D-O5D-PN-O3
2	D	501	NDP	C5D-O5D-PN-O1N
2	C	501	NDP	C5B-O5B-PA-O1A
2	C	501	NDP	C5B-O5B-PA-O2A
2	C	501	NDP	C5B-O5B-PA-O3
2	C	501	NDP	C5D-O5D-PN-O3
2	C	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	O4D-C1D-N1N-C6N
2	C	501	NDP	C3D-C4D-C5D-O5D
2	D	501	NDP	O4D-C1D-N1N-C2N
2	D	501	NDP	C2D-C1D-N1N-C6N
2	A	501	NDP	C3D-C4D-C5D-O5D
2	C	501	NDP	O4D-C4D-C5D-O5D
2	D	501	NDP	C2D-C1D-N1N-C2N
2	A	501	NDP	C3B-C4B-C5B-O5B
2	A	501	NDP	O4D-C4D-C5D-O5D
2	B	501	NDP	O4B-C4B-C5B-O5B
2	C	501	NDP	C4D-C5D-O5D-PN
2	A	501	NDP	PN-O3-PA-O5B
2	A	501	NDP	PA-O3-PN-O5D
2	C	501	NDP	PA-O3-PN-O5D
2	A	501	NDP	O4B-C4B-C5B-O5B
2	D	501	NDP	C4D-C5D-O5D-PN
2	D	501	NDP	C5B-O5B-PA-O1A
2	C	501	NDP	C5D-O5D-PN-O1N
2	C	501	NDP	C2N-C3N-C7N-N7N
2	C	501	NDP	C2B-O2B-P2B-O1X
2	A	501	NDP	PN-O3-PA-O1A
2	C	501	NDP	PA-O3-PN-O1N
2	C	501	NDP	C2B-O2B-P2B-O2X
2	B	501	NDP	C2N-C3N-C7N-O7N

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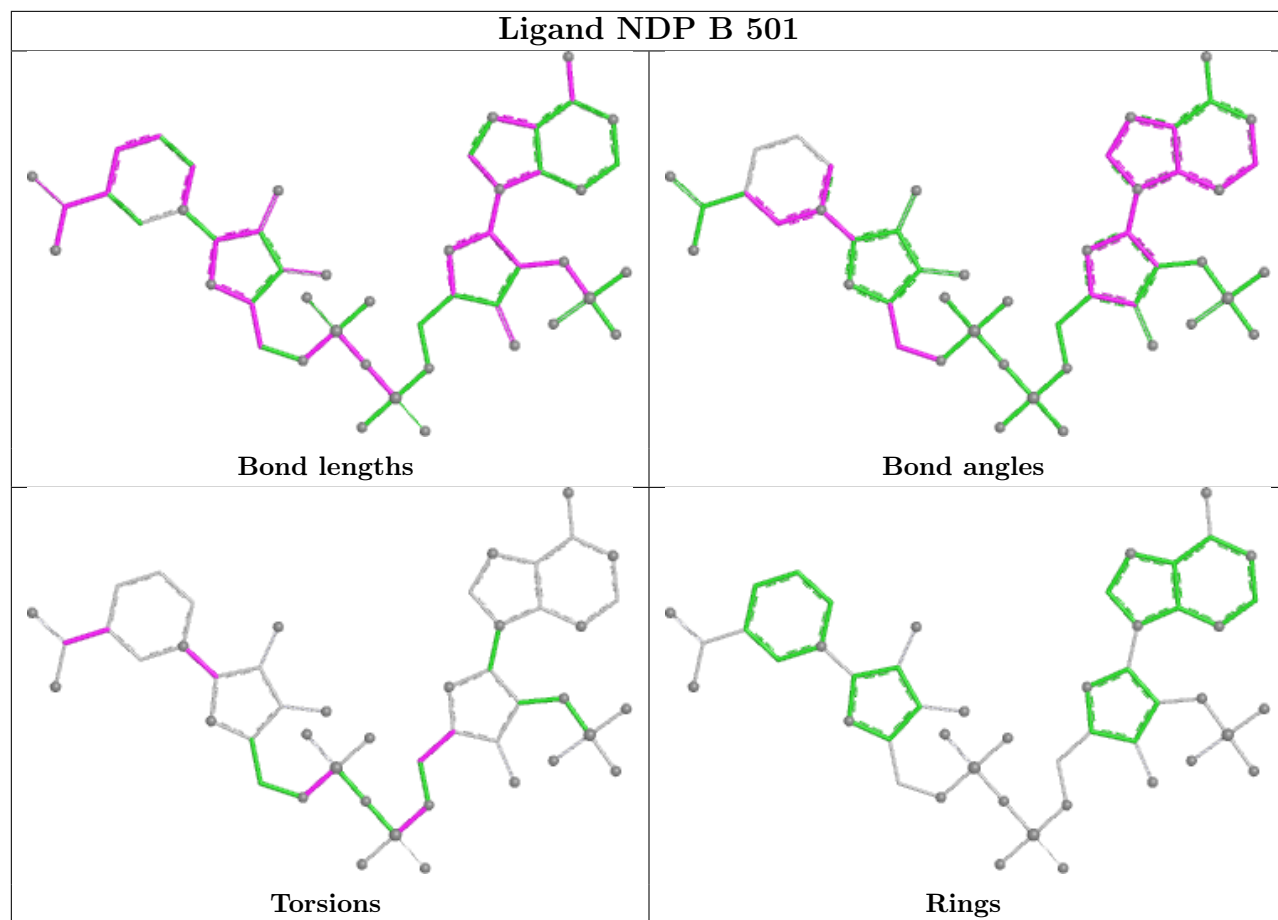
Mol	Chain	Res	Type	Atoms
2	D	501	NDP	O4B-C1B-N9A-C8A
2	C	501	NDP	O4B-C4B-C5B-O5B

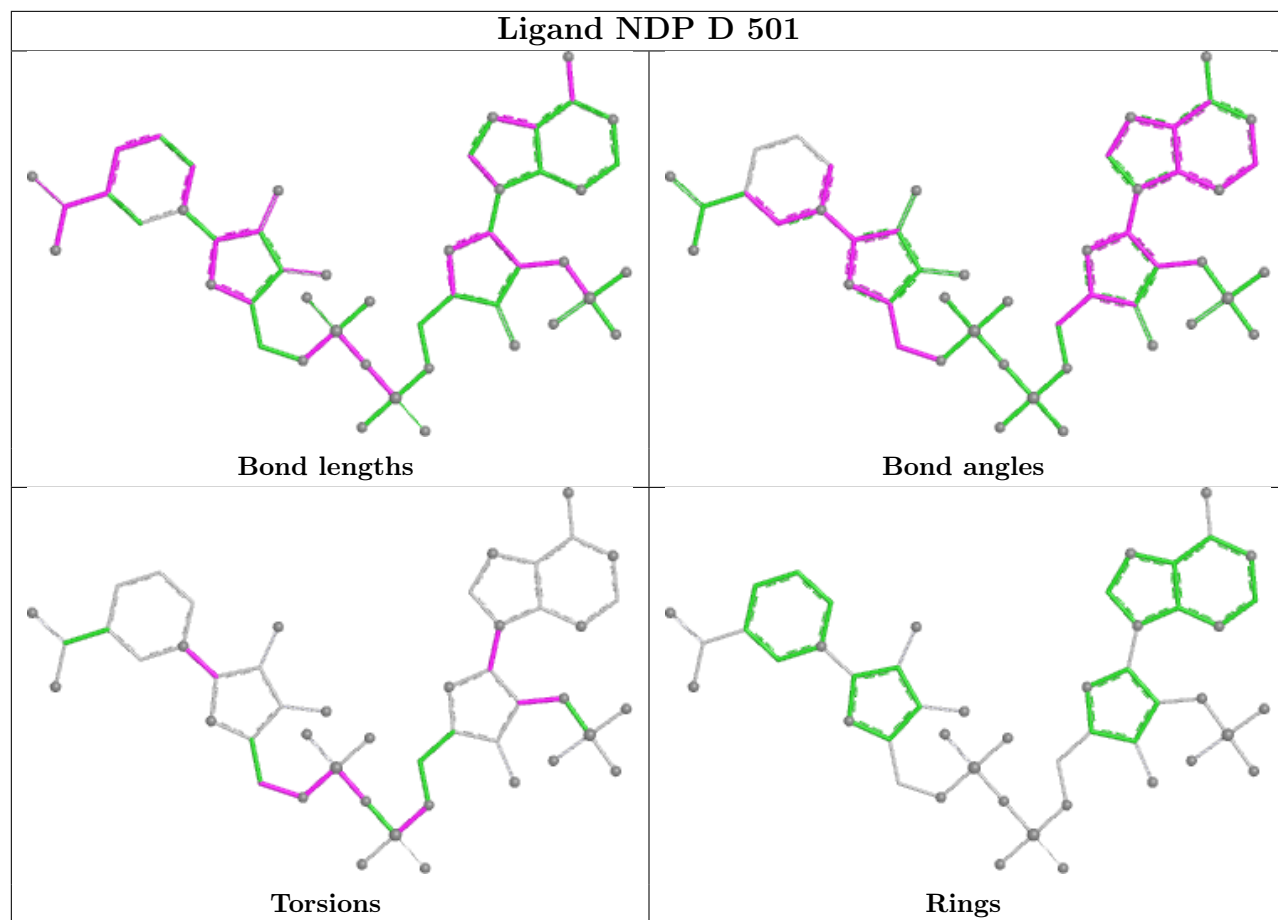
There are no ring outliers.

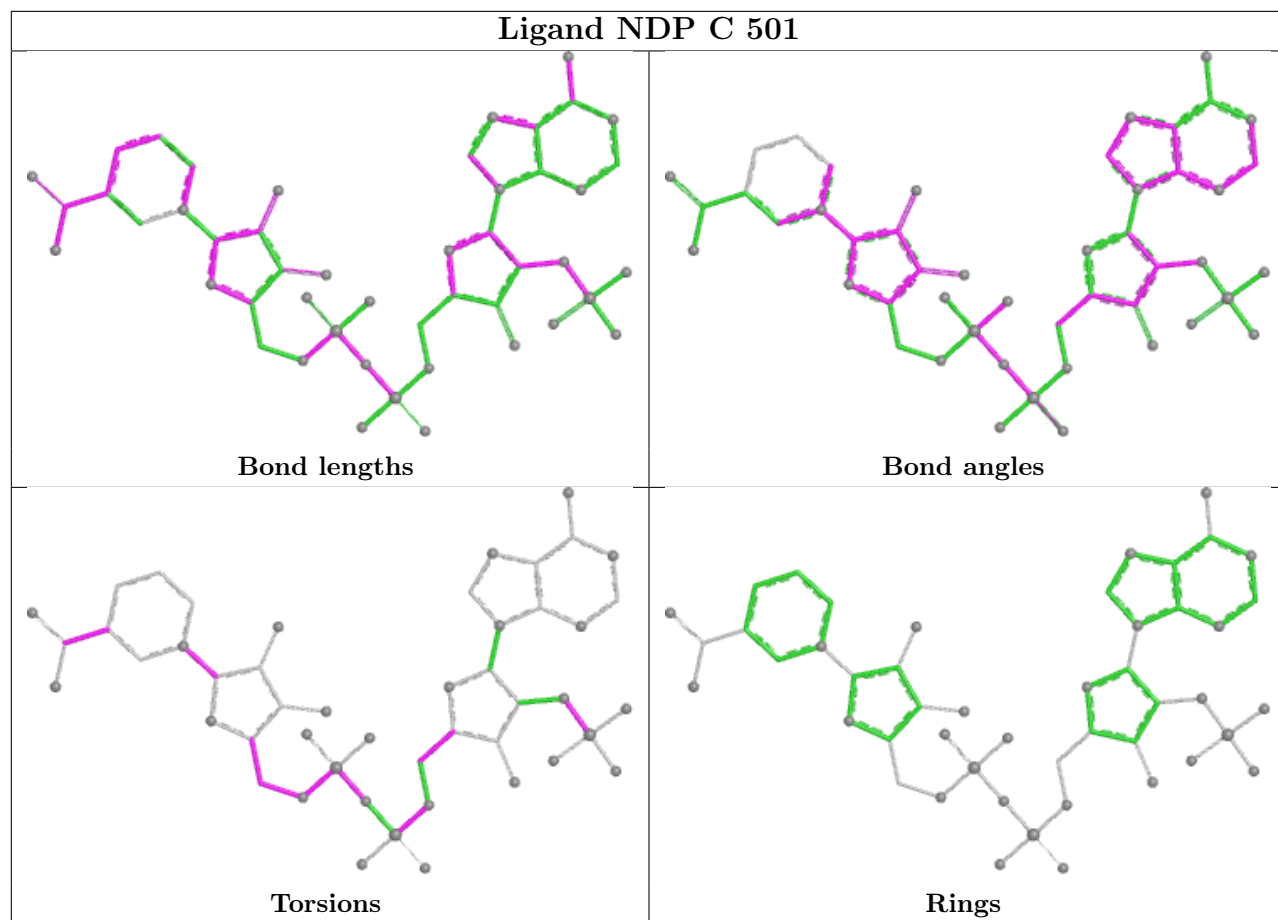
4 monomers are involved in 35 short contacts:

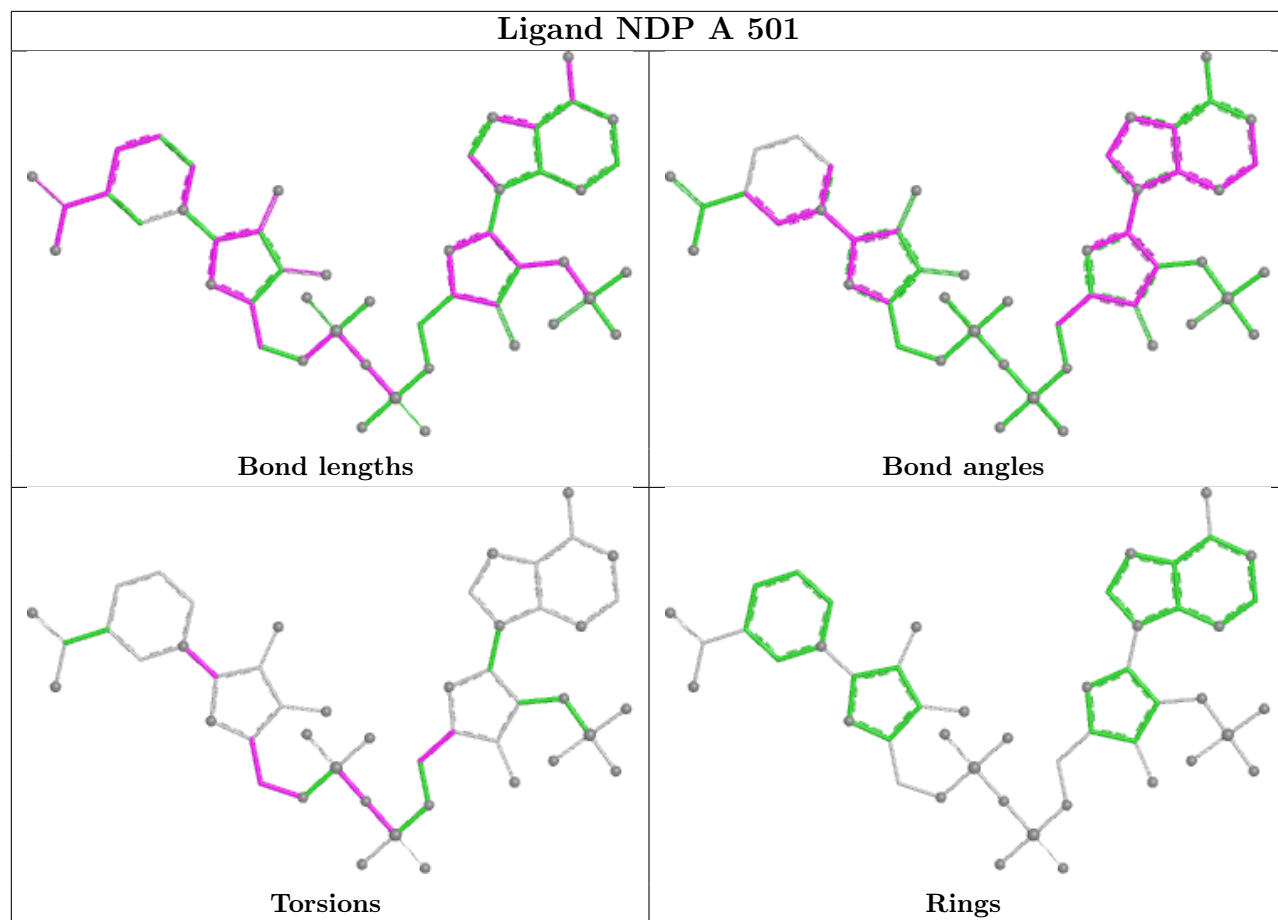
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	NDP	3	0
2	D	501	NDP	16	0
2	C	501	NDP	5	0
2	A	501	NDP	11	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	445/445 (100%)	1.02	73 (16%) 4 4	10, 36, 62, 80	1 (0%)
1	B	444/445 (99%)	0.80	43 (9%) 13 14	13, 30, 54, 79	1 (0%)
1	C	445/445 (100%)	0.85	47 (10%) 11 12	13, 31, 53, 72	2 (0%)
1	D	444/445 (99%)	1.22	84 (18%) 3 3	10, 37, 74, 110	1 (0%)
All	All	1778/1780 (99%)	0.97	247 (13%) 6 6	10, 33, 62, 110	5 (0%)

All (247) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	103	LEU	8.4
1	D	103	LEU	7.7
1	D	96	PHE	7.6
1	D	43	LEU	7.3
1	A	15	SER	6.6
1	D	40	PHE	6.6
1	D	97	LEU	5.5
1	D	166	ARG	5.4
1	D	95	GLY	5.4
1	C	40	PHE	5.3
1	D	42	GLY	5.3
1	D	44	ALA	5.1
1	D	112	LEU	5.1
1	D	169	GLY	5.0
1	A	44	ALA	5.0
1	D	100	TYR	5.0
1	B	87	ILE	4.8
1	B	43	LEU	4.8
1	B	95	GLY	4.6
1	D	167	ILE	4.4
1	D	84	TRP	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	339	CYS	4.3
1	D	39	MET	4.2
1	D	114	TYR	4.2
1	D	45	SER	4.1
1	D	46	ARG	4.0
1	A	101	GLY	4.0
1	C	409	HIS	3.9
1	A	13	ALA	3.9
1	A	80	TYR	3.9
1	C	100	TYR	3.9
1	D	80	TYR	3.8
1	D	419	ALA	3.8
1	D	106	LEU	3.6
1	A	20	ALA	3.6
1	B	46	ARG	3.6
1	B	100	TYR	3.6
1	D	99	ARG	3.6
1	A	132	VAL	3.5
1	D	22	ALA	3.5
1	C	43	LEU	3.5
1	C	53[A]	SER	3.5
1	D	102	ARG	3.5
1	B	94	PHE	3.4
1	D	156	HIS	3.3
1	D	48	LYS	3.3
1	A	11	GLY	3.3
1	D	94	PHE	3.3
1	D	41	ASP	3.3
1	A	87	ILE	3.3
1	D	366	PHE	3.3
1	D	12	ASP	3.3
1	D	50	PRO	3.3
1	A	433	SER	3.2
1	C	35	GLY	3.2
1	A	90	PRO	3.2
1	D	98	GLU	3.2
1	D	110	HIS	3.2
1	A	424	LEU	3.1
1	A	172	THR	3.1
1	C	410	GLN	3.1
1	C	256	VAL	3.1
1	B	40	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	93	THR	3.1
1	C	23	LEU	3.1
1	C	411	GLY	3.1
1	C	116	VAL	3.1
1	A	56	LEU	3.1
1	B	96	PHE	3.0
1	A	29	ALA	3.0
1	B	161	MET	3.0
1	D	92	SER	3.0
1	C	355	TRP	3.0
1	D	101	GLY	3.0
1	A	88	PHE	3.0
1	C	96	PHE	3.0
1	D	16	ALA	3.0
1	D	267	ALA	3.0
1	C	29	ALA	3.0
1	A	43	LEU	3.0
1	A	42	GLY	3.0
1	D	264	ILE	3.0
1	B	93	THR	3.0
1	C	114	TYR	3.0
1	B	39	MET	3.0
1	A	404	VAL	2.9
1	C	264	ILE	2.9
1	B	47	ASP	2.9
1	D	1	MET	2.9
1	C	41	ASP	2.9
1	A	399	GLN	2.9
1	B	88	PHE	2.9
1	A	22	ALA	2.9
1	D	19	ALA	2.9
1	A	97	LEU	2.9
1	B	132	VAL	2.9
1	A	161	MET	2.9
1	A	40	PHE	2.9
1	B	84	TRP	2.9
1	C	102	ARG	2.8
1	D	429	PRO	2.8
1	D	104	SER	2.8
1	A	388	LEU	2.8
1	A	23	LEU	2.8
1	A	30	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	404	VAL	2.8
1	A	293	ARG	2.8
1	D	165	GLN	2.8
1	C	262	ALA	2.8
1	A	237	LEU	2.8
1	D	436	LEU	2.8
1	D	107	THR	2.8
1	D	132	VAL	2.8
1	D	13	ALA	2.7
1	D	178	ALA	2.7
1	A	83	VAL	2.7
1	A	405	HIS	2.7
1	D	91	VAL	2.7
1	D	115	HIS	2.7
1	C	234	SER	2.7
1	C	101	GLY	2.7
1	A	21	LEU	2.7
1	D	182	LEU	2.7
1	B	105	PRO	2.7
1	A	28	ARG	2.7
1	B	78	VAL	2.6
1	C	407	ASN	2.6
1	D	52	LYS	2.6
1	D	369	TYR	2.6
1	C	22	ALA	2.6
1	D	59	VAL	2.6
1	A	123	GLY	2.6
1	B	13	ALA	2.6
1	D	177	LEU	2.6
1	B	442	PHE	2.5
1	A	8	ILE	2.5
1	D	176	GLY	2.5
1	B	443	ARG	2.5
1	D	20	ALA	2.5
1	A	402	LEU	2.5
1	D	88	PHE	2.5
1	A	141	VAL	2.5
1	A	415	VAL	2.5
1	D	83	VAL	2.5
1	C	288	HIS	2.5
1	D	47	ASP	2.5
1	A	62	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	90	PRO	2.5
1	B	101	GLY	2.5
1	A	255	VAL	2.5
1	C	166	ARG	2.5
1	A	14	ALA	2.5
1	A	420	PRO	2.5
1	A	436	LEU	2.5
1	A	59	VAL	2.4
1	A	12	ASP	2.4
1	B	58	ASP	2.4
1	B	80	TYR	2.4
1	A	262	ALA	2.4
1	A	61	LEU	2.4
1	B	398	GLY	2.4
1	B	114	TYR	2.4
1	A	406	HIS	2.4
1	D	111	ASP	2.4
1	D	234	SER	2.4
1	D	143	ALA	2.4
1	A	411	GLY	2.4
1	A	174	PHE	2.4
1	A	102	ARG	2.4
1	C	99	ARG	2.4
1	B	254	CYS	2.4
1	C	84	TRP	2.4
1	A	41	ASP	2.3
1	C	434	LYS	2.3
1	B	90	PRO	2.3
1	A	103	LEU	2.3
1	A	126	LEU	2.3
1	C	131	GLY	2.3
1	D	125	VAL	2.3
1	B	9	LEU	2.3
1	C	197	TRP	2.3
1	A	45	SER	2.3
1	C	0	HIS	2.3
1	D	71	VAL	2.3
1	A	24	PRO	2.3
1	A	106	LEU	2.3
1	C	168	TRP	2.3
1	B	305	PHE	2.2
1	B	8	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	58	ASP	2.2
1	D	14	ALA	2.2
1	D	200	ALA	2.2
1	C	278	ALA	2.2
1	A	168	TRP	2.2
1	A	104	SER	2.2
1	C	303	ARG	2.2
1	D	439	ILE	2.2
1	B	329	VAL	2.2
1	B	42	GLY	2.2
1	B	98	GLU	2.2
1	C	431	LEU	2.2
1	B	45	SER	2.2
1	B	197	TRP	2.2
1	D	320	ARG	2.2
1	C	-1	ARG	2.2
1	B	12	ASP	2.2
1	C	160	MET	2.2
1	C	404	VAL	2.2
1	D	35	GLY	2.2
1	D	150	LEU	2.2
1	D	38	ARG	2.2
1	D	62	PRO	2.1
1	D	37	GLU	2.1
1	C	91	VAL	2.1
1	A	425	GLY	2.1
1	A	400	ALA	2.1
1	D	130	ALA	2.1
1	C	7	ALA	2.1
1	C	20	ALA	2.1
1	A	146	LEU	2.1
1	C	222	LEU	2.1
1	B	26	SER	2.1
1	C	104	SER	2.1
1	A	182	LEU	2.1
1	A	173	ASN	2.1
1	C	292	PRO	2.1
1	A	155	GLY	2.1
1	B	273	ILE	2.1
1	A	96	PHE	2.1
1	B	32	LEU	2.1
1	B	36	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	46	ARG	2.1
1	B	102	ARG	2.1
1	A	36	GLU	2.0
1	B	173	ASN	2.0
1	A	409	HIS	2.0
1	A	131	GLY	2.0
1	C	392	TYR	2.0
1	A	391	VAL	2.0
1	D	311	VAL	2.0
1	C	413	VAL	2.0
1	D	168	TRP	2.0
1	D	284	TRP	2.0
1	A	5	LEU	2.0
1	D	146	LEU	2.0
1	A	107	THR	2.0
1	D	105	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

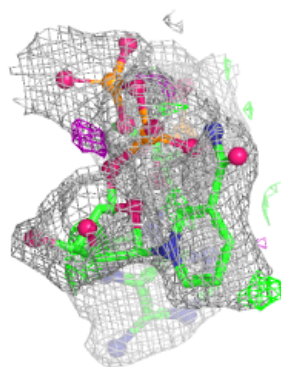
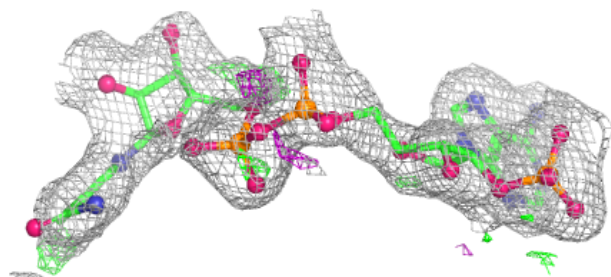
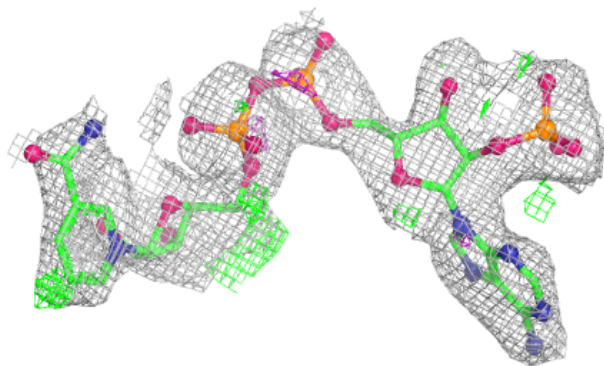
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NDP	A	501	48/48	0.86	0.15	29,53,69,85	0
2	NDP	C	501	48/48	0.89	0.11	12,28,42,55	0
2	NDP	D	501	48/48	0.90	0.14	24,42,60,62	0
2	NDP	B	501	48/48	0.93	0.11	15,35,70,76	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

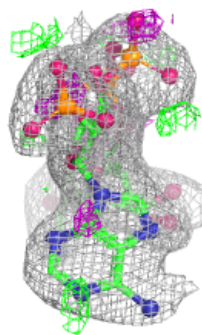
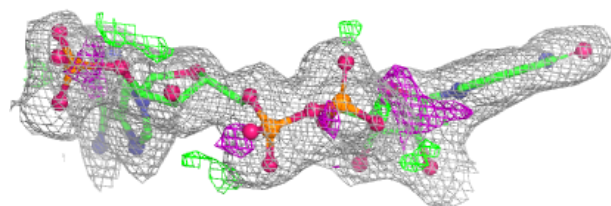
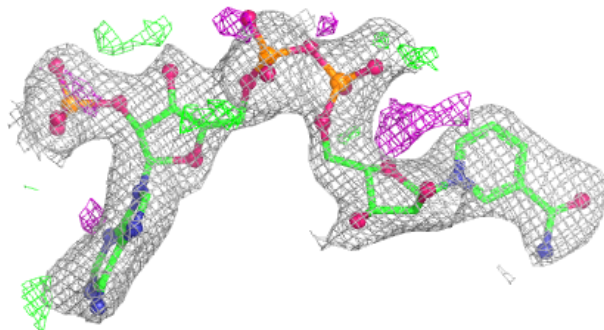
Electron density around NDP A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



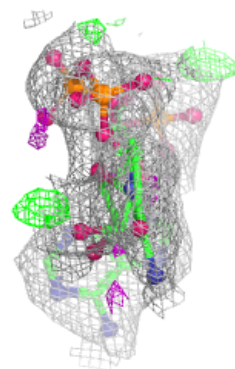
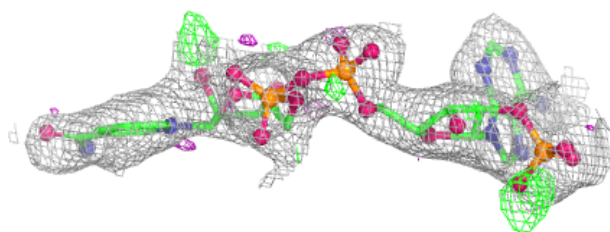
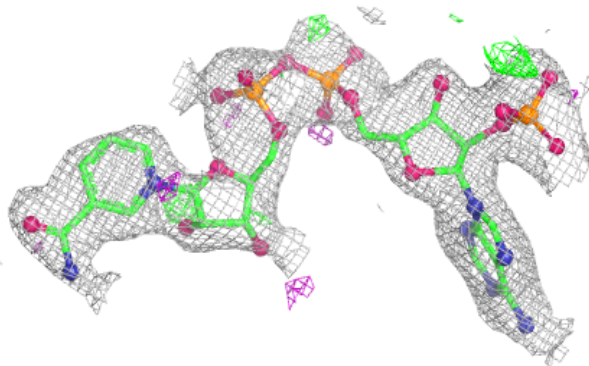
Electron density around NDP C 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



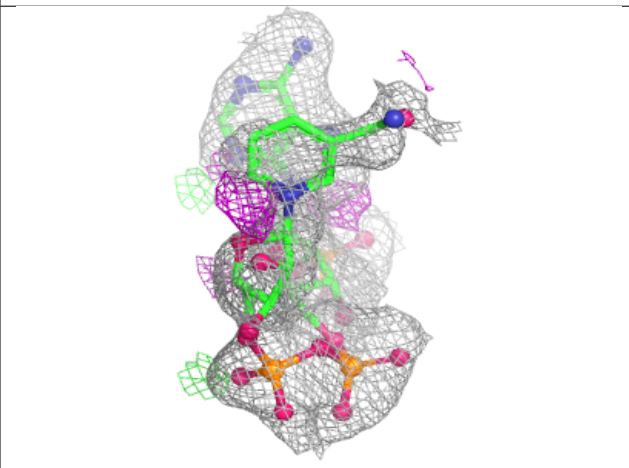
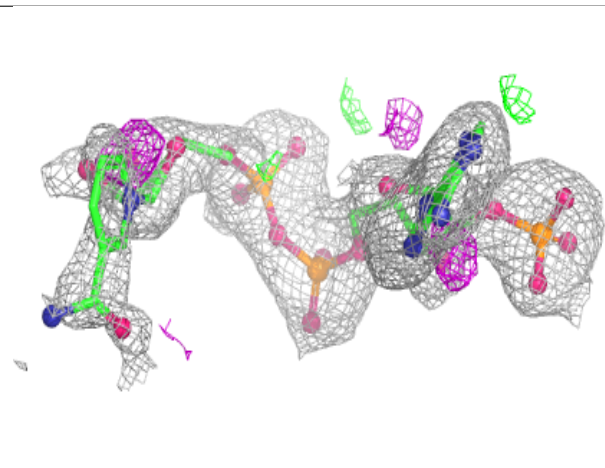
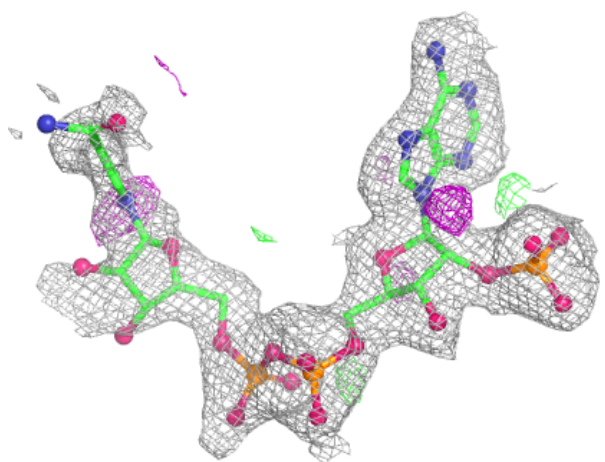
Electron density around NDP D 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NDP B 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.