



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:50 AM UTC

PDB ID : 4CJC / pdb_00004cjc
Title : orthorhombic crystal form of Bogt6a E192Q in complex with UDP-GalNAc, UDP, GalNAc
Authors : Pham, T.; Stinson, B.; Thiyagarajan, N.; Lizotte-Waniewski, M.; Brew, K.; Acharya, K.R.
Deposited on : 2013-12-19
Resolution : 3.42 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

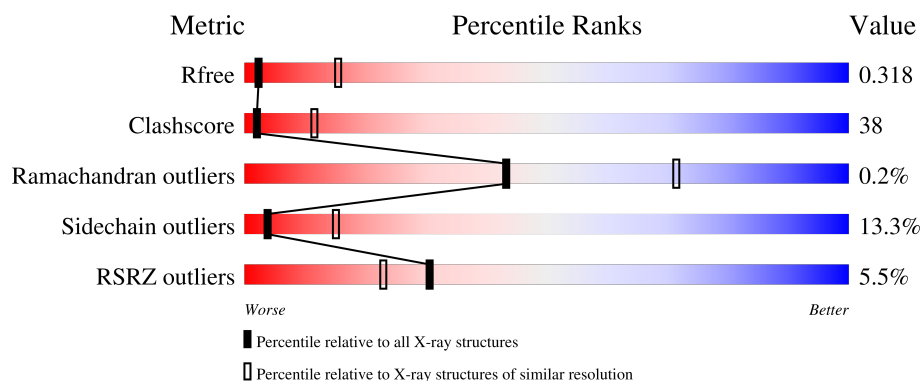
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.42 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	1210 (3.48-3.36)
Clashscore	190562	1234 (3.48-3.36)
Ramachandran outliers	187476	1222 (3.48-3.36)
Sidechain outliers	187428	1222 (3.48-3.36)
RSRZ outliers	180081	1210 (3.48-3.36)

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	246	0% 47% 38% 12% .
1	B	246	2% 50% 39% 8% .
1	C	246	3% 49% 36% 12% ..
1	D	246	16% 39% 50% 7% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8120 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.

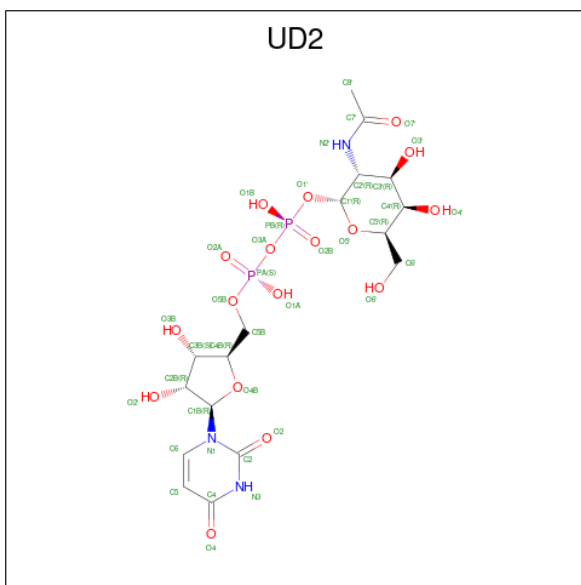
- Molecule 1 is a protein called GLYCOSYLTRANSFERASE FAMILY 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	7	0	0
			2008	1312	327	362	7			
1	B	238	Total	C	N	O	S	10	0	0
			1981	1295	322	357	7			
1	C	240	Total	C	N	O	S	10	0	0
			2000	1306	326	361	7			
1	D	236	Total	C	N	O	S	10	0	0
			1973	1291	320	355	7			

There are 4 discrepancies between the modelled and reference sequences:

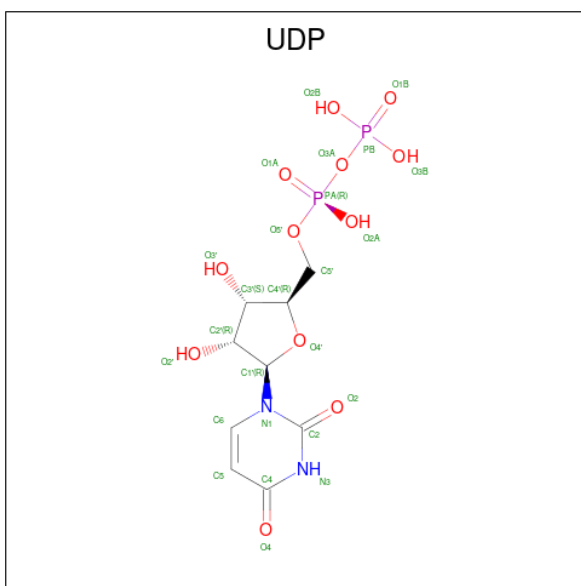
Chain	Residue	Modelled	Actual	Comment	Reference
A	192	GLN	GLU	engineered mutation	UNP A7LVT2
B	192	GLN	GLU	engineered mutation	UNP A7LVT2
C	192	GLN	GLU	engineered mutation	UNP A7LVT2
D	192	GLN	GLU	engineered mutation	UNP A7LVT2

- Molecule 2 is URIDINE-DIPHOSPHATE-N-ACETYLGALACTOSAMINE (CCD ID: UD2) (formula: C₁₇H₂₇N₃O₁₇P₂).



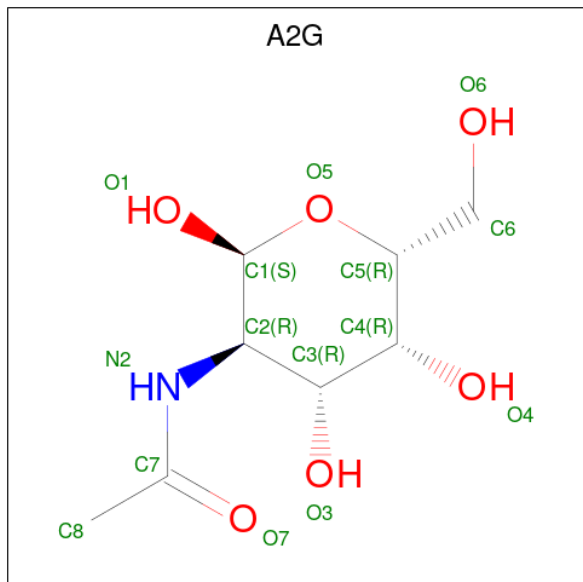
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 39	C 17	N 3	O 17	P 2	0	0
2	C	1	Total 39	C 17	N 3	O 17	P 2	0	0

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: $\text{C}_9\text{H}_{14}\text{N}_2\text{O}_{12}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total 25	C 9	N 2	O 12	P 2	0	0
3	D	1	Total 25	C 9	N 2	O 12	P 2	0	0

- Molecule 4 is 2-acetamido-2-deoxy- α -D-galactopyranose (CCD ID: A2G) (formula: $C_8H_{15}NO_6$).

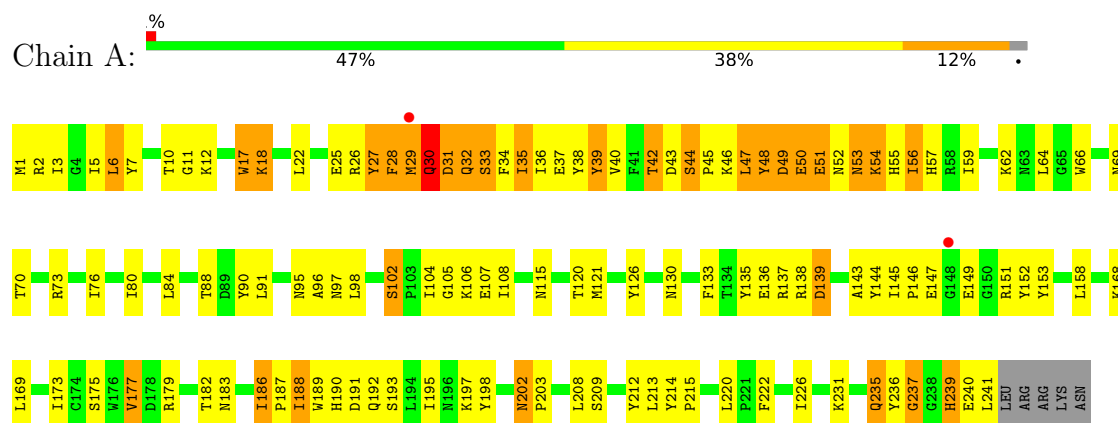


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			15	8	1	6		
4	D	1	Total	C	N	O	0	0
			15	8	1	6		

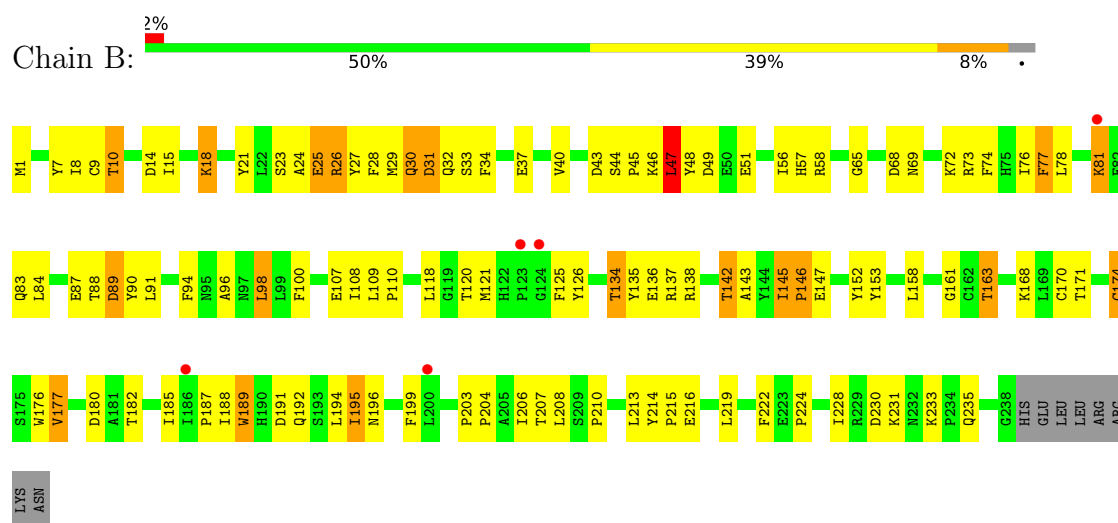
3 Residue-property plots

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.

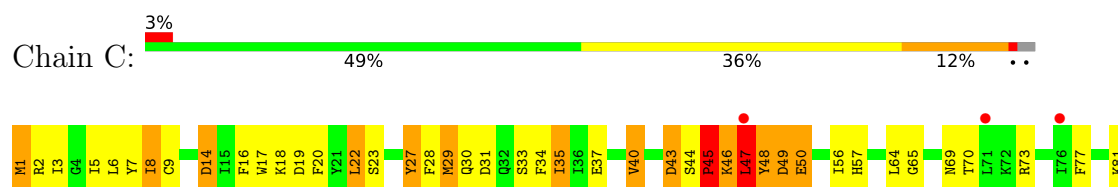
• Molecule 1: GLYCOSYLTRANSFERASE FAMILY 6

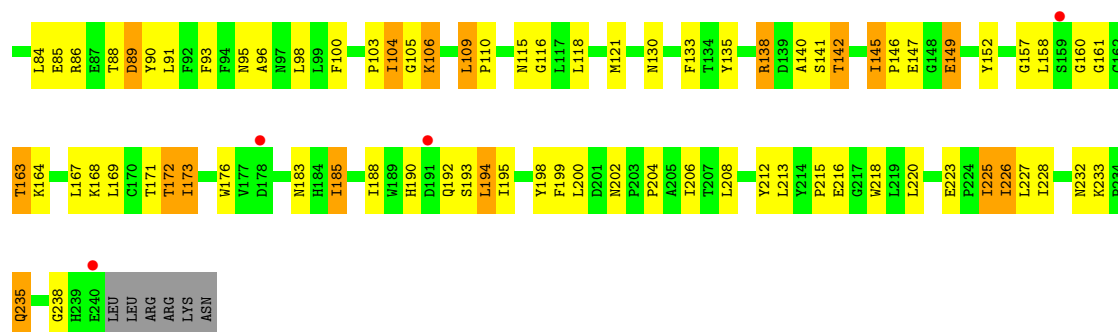


• Molecule 1: GLYCOSYLTRANSFERASE FAMILY 6

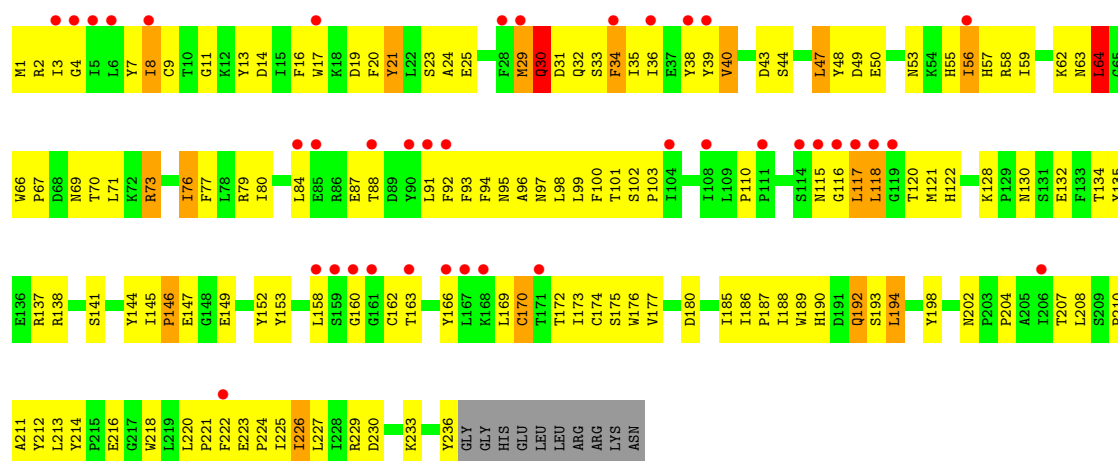


• Molecule 1: GLYCOSYLTRANSFERASE FAMILY 6





• Molecule 1: GLYCOSYLTRANSFERASE FAMILY 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.12Å 120.15Å 131.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.80 – 3.42 88.80 – 3.42	Depositor EDS
% Data completeness (in resolution range)	97.0 (88.80-3.42) 90.4 (88.80-3.42)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 3.41Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.283 , 0.314 0.292 , 0.318	Depositor DCC
R_{free} test set	878 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	84.6	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	8120	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UD2, UDP, A2G

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	0.67	0/2074	0.96	11/2816 (0.4%)
1	B	0.58	0/2046	0.96	10/2778 (0.4%)
1	C	0.45	0/2066	0.92	5/2805 (0.2%)
1	D	0.44	0/2038	0.99	5/2768 (0.2%)
All	All	0.54	0/8224	0.96	31/11167 (0.3%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49	ASP	N-CA-C	-7.93	88.78	111.00
1	A	239	HIS	N-CA-C	7.68	121.34	108.20
1	B	77	PHE	N-CA-C	-7.57	103.93	113.01
1	B	34	PHE	N-CA-CB	7.30	121.39	110.29
1	A	48	TYR	N-CA-C	7.22	120.69	110.23
1	D	73	ARG	N-CA-C	-7.10	103.54	111.71
1	D	170	CYS	N-CA-C	6.97	118.88	111.28
1	A	237	GLY	N-CA-C	-6.74	97.20	113.18
1	A	30	GLN	N-CA-C	6.61	118.48	111.28
1	A	31	ASP	N-CA-C	-6.55	99.62	110.17
1	A	17	TRP	N-CA-C	6.54	118.41	111.28
1	C	85	GLU	N-CA-C	-6.44	105.55	113.41
1	B	30	GLN	N-CA-C	6.20	120.16	112.59
1	A	186	ILE	CB-CA-C	-5.99	104.32	110.13
1	D	30	GLN	N-CA-C	5.95	117.77	111.28
1	B	33	SER	CA-C-O	5.92	126.74	119.05
1	B	30	GLN	CB-CA-C	5.91	120.54	111.02
1	C	47	LEU	CA-C-N	-5.83	114.43	122.30
1	C	47	LEU	C-N-CA	-5.83	114.43	122.30
1	C	27	TYR	N-CA-C	5.81	121.80	114.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	45	PRO	CB-CA-C	-5.76	103.91	112.55
1	B	33	SER	CB-CA-C	-5.65	98.64	110.17
1	A	188	ILE	CB-CA-C	-5.63	104.76	111.97
1	B	32	GLN	N-CA-C	5.63	117.86	111.11
1	A	18	LYS	N-CA-C	5.43	116.89	110.97
1	D	64	LEU	N-CA-C	-5.37	106.74	113.72
1	B	26	ARG	N-CA-C	5.31	117.15	111.36
1	B	31	ASP	N-CA-C	-5.29	102.56	110.23
1	A	27	TYR	N-CA-C	5.28	120.21	113.55
1	D	34	PHE	N-CA-C	5.25	118.14	109.85
1	B	47	LEU	CA-CB-CG	-5.04	98.66	116.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2008	0	1944	200	0
1	B	1981	0	1920	103	0
1	C	2000	0	1933	136	0
1	D	1973	0	1914	162	0
2	A	39	0	25	9	0
2	C	39	0	25	5	0
3	B	25	0	11	2	0
3	D	25	0	11	2	0
4	B	15	0	12	0	0
4	D	15	0	12	2	0
All	All	8120	0	7807	600	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 38.

All (600) 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:D:29:MET:HG3	1:D:34:PHE:CZ	1.33	1.63
1:C:48:TYR:CD1	1:C:49:ASP:HB3	1.41	1.52
1:A:3:ILE:CD1	1:A:29:MET:HE3	1.42	1.48
1:A:3:ILE:HD12	1:A:29:MET:CE	1.41	1.46
1:A:188:ILE:HD13	1:A:189:TRP:CD1	1.49	1.44
1:A:3:ILE:CD1	1:A:29:MET:CE	1.95	1.43
1:A:3:ILE:HD11	1:A:29:MET:SD	1.61	1.40
1:D:29:MET:HG3	1:D:34:PHE:CE2	1.55	1.38
1:B:28:PHE:CE1	1:B:29:MET:HG3	1.61	1.35
1:C:48:TYR:CE1	1:C:49:ASP:HB3	1.62	1.32
1:A:188:ILE:HD12	1:A:189:TRP:N	1.48	1.26
1:C:29:MET:O	1:C:34:PHE:HZ	1.14	1.24
1:D:29:MET:CG	1:D:34:PHE:CE2	2.18	1.24
1:A:50:GLU:O	1:A:51:GLU:HG2	1.34	1.22
1:C:29:MET:O	1:C:34:PHE:CZ	1.92	1.22
1:A:188:ILE:HD12	1:A:189:TRP:CB	1.69	1.20
1:D:29:MET:HB3	1:D:34:PHE:CE2	1.76	1.19
1:D:29:MET:CG	1:D:34:PHE:CZ	2.26	1.18
1:A:29:MET:HB2	1:A:36:ILE:HD11	1.26	1.16
1:A:42:THR:HG23	1:A:44:SER:H	1.10	1.15
1:A:188:ILE:CD1	1:A:189:TRP:CD1	2.30	1.15
1:C:48:TYR:CE1	1:C:49:ASP:CB	2.30	1.15
1:D:50:GLU:OE1	1:D:58:ARG:NH2	1.78	1.14
1:C:28:PHE:O	1:C:29:MET:HG2	1.47	1.14
1:A:49:ASP:OD1	1:A:49:ASP:O	1.66	1.14
1:B:28:PHE:CE1	1:B:29:MET:CG	2.30	1.14
1:D:29:MET:CB	1:D:34:PHE:CE2	2.32	1.13
1:A:54:LYS:HD2	1:A:54:LYS:O	1.50	1.12
1:A:188:ILE:CD1	1:A:189:TRP:HB2	1.78	1.12
1:A:3:ILE:CD1	1:A:29:MET:SD	2.32	1.11
1:A:26:ARG:HD3	1:A:27:TYR:HE1	1.14	1.10
1:B:46:LYS:O	1:B:47:LEU:HD23	1.48	1.10
1:D:29:MET:HB3	1:D:34:PHE:CD2	1.86	1.09
1:C:48:TYR:CD1	1:C:48:TYR:C	2.30	1.09
1:A:188:ILE:CD1	1:A:189:TRP:CB	2.30	1.08
1:A:26:ARG:HD3	1:A:27:TYR:CE1	1.88	1.08
1:C:47:LEU:HD22	1:C:56:ILE:HG21	1.31	1.08
1:C:48:TYR:CD1	1:C:49:ASP:CB	2.37	1.07
1:A:188:ILE:HD12	1:A:189:TRP:CA	1.84	1.06
1:D:40:VAL:HG11	1:D:47:LEU:HD21	1.31	1.06
1:A:40:VAL:HG11	1:A:47:LEU:HD11	1.09	1.06
1:D:40:VAL:HG11	1:D:47:LEU:CD2	1.86	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:GLU:CD	1:A:51:GLU:H	1.68	1.02
1:C:48:TYR:HD1	1:C:48:TYR:O	1.41	1.02
1:A:17:TRP:NE1	1:A:48:TYR:O	1.93	1.00
1:B:26:ARG:NH1	1:B:27:TYR:CZ	2.30	1.00
1:D:47:LEU:N	1:D:58:ARG:HH21	1.58	1.00
1:A:29:MET:HG2	1:A:36:ILE:HG12	1.41	1.00
1:C:45:PRO:HG2	1:C:46:LYS:N	1.76	0.99
1:A:28:PHE:CE2	1:A:29:MET:HE2	1.98	0.98
1:A:48:TYR:CE1	1:A:49:ASP:HB3	1.99	0.97
1:A:40:VAL:CG1	1:A:47:LEU:HD11	1.94	0.96
1:A:188:ILE:CD1	1:A:189:TRP:CG	2.46	0.96
1:D:47:LEU:H	1:D:58:ARG:NH2	1.62	0.96
1:D:120:THR:HG22	1:D:208:LEU:HB2	1.43	0.96
1:C:45:PRO:CD	1:C:46:LYS:H	1.79	0.96
1:A:28:PHE:CD2	1:A:29:MET:HE2	2.00	0.94
1:C:48:TYR:C	1:C:48:TYR:HD1	1.71	0.94
1:A:42:THR:HG23	1:A:44:SER:N	1.81	0.94
1:A:188:ILE:CD1	1:A:189:TRP:N	2.30	0.94
1:C:45:PRO:CG	1:C:46:LYS:N	2.30	0.94
1:B:28:PHE:CD1	1:B:29:MET:CG	2.51	0.94
1:A:3:ILE:HD12	1:A:29:MET:HE1	1.48	0.93
1:A:188:ILE:HD11	1:A:189:TRP:HB2	1.51	0.93
1:B:28:PHE:HE1	1:B:29:MET:HG3	1.01	0.92
1:A:3:ILE:HD12	1:A:29:MET:HE3	0.95	0.92
1:C:45:PRO:CG	1:C:46:LYS:H	1.82	0.92
1:D:13:TYR:HD2	3:D:1237:UDP:H2'	1.35	0.92
1:A:42:THR:CG2	1:A:44:SER:N	2.32	0.91
1:A:42:THR:CG2	1:A:44:SER:H	1.83	0.91
1:A:50:GLU:CD	1:A:51:GLU:N	2.28	0.91
1:A:29:MET:CB	1:A:36:ILE:HD11	2.01	0.90
1:A:25:GLU:OE2	1:A:38:TYR:OH	1.88	0.89
1:C:47:LEU:CD2	1:C:56:ILE:HG21	2.03	0.89
1:B:28:PHE:HD1	1:B:29:MET:HB2	1.37	0.88
1:A:53:ASN:OD1	1:A:56:ILE:HG22	1.74	0.88
1:D:47:LEU:N	1:D:58:ARG:NH2	2.22	0.88
1:A:10:THR:H	2:A:1242:UD2:H5	1.35	0.87
1:A:29:MET:HB2	1:A:36:ILE:CD1	2.02	0.87
1:C:20:PHE:HE1	1:C:100:PHE:HE1	1.23	0.87
1:A:3:ILE:HD11	1:A:29:MET:CE	1.80	0.86
1:C:48:TYR:HD1	1:C:49:ASP:HB3	1.14	0.86
1:A:188:ILE:HD12	1:A:189:TRP:CG	2.08	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:VAL:HG11	1:A:47:LEU:CD1	2.03	0.85
1:A:29:MET:CG	1:A:36:ILE:CG1	2.55	0.84
1:A:50:GLU:HA	1:A:50:GLU:OE1	1.74	0.84
1:C:47:LEU:HB3	1:C:50:GLU:OE2	1.78	0.84
1:C:48:TYR:HE1	1:C:49:ASP:CB	1.88	0.83
1:A:37:GLU:OE2	1:A:55:HIS:ND1	2.10	0.83
1:A:50:GLU:C	1:A:51:GLU:HG2	2.02	0.83
1:C:29:MET:C	1:C:34:PHE:CZ	2.56	0.82
1:A:188:ILE:HD13	1:A:189:TRP:HD1	0.98	0.82
1:D:1:MET:N	1:D:33:SER:O	2.11	0.82
1:A:192:GLN:HE21	2:A:1242:UD2:H3'	1.44	0.81
1:A:26:ARG:HB3	1:A:27:TYR:CD1	2.16	0.81
1:A:50:GLU:OE1	1:A:50:GLU:CA	2.29	0.81
1:D:47:LEU:HB3	1:D:56:ILE:HD11	1.63	0.81
1:A:3:ILE:HD11	1:A:29:MET:CG	2.11	0.81
1:D:19:ASP:HB2	1:D:229:ARG:HH21	1.46	0.81
1:A:231:LYS:NZ	2:A:1242:UD2:O3A	2.15	0.80
1:D:84:LEU:O	1:D:88:THR:OG1	2.00	0.80
1:A:179:ARG:O	1:A:183:ASN:ND2	2.15	0.80
1:C:47:LEU:CB	1:C:50:GLU:OE2	2.29	0.80
1:A:29:MET:CG	1:A:36:ILE:HG12	2.12	0.79
1:D:13:TYR:CD2	3:D:1237:UDP:H2'	2.17	0.79
1:A:54:LYS:O	1:A:54:LYS:CD	2.30	0.79
1:B:46:LYS:O	1:B:47:LEU:CD2	2.30	0.79
1:A:17:TRP:CD1	1:A:47:LEU:HD23	2.16	0.79
1:C:45:PRO:HD2	1:C:46:LYS:H	1.45	0.79
1:C:29:MET:C	1:C:34:PHE:HZ	1.90	0.79
1:C:47:LEU:HB3	1:C:50:GLU:CD	2.09	0.78
1:A:29:MET:HG2	1:A:36:ILE:CG1	2.13	0.78
1:B:84:LEU:HD23	1:B:91:LEU:HD11	1.65	0.78
1:D:17:TRP:HE1	1:D:48:TYR:HD1	1.30	0.78
1:B:28:PHE:CD1	1:B:29:MET:HB2	2.19	0.77
1:B:121:MET:HB2	1:B:210:PRO:HD3	1.66	0.77
1:A:53:ASN:OD1	1:A:56:ILE:CG2	2.32	0.77
1:C:45:PRO:HG2	1:C:46:LYS:H	1.39	0.77
1:A:48:TYR:CD1	1:A:49:ASP:HB3	2.19	0.77
1:A:54:LYS:HD2	1:A:54:LYS:C	2.10	0.77
1:D:50:GLU:OE2	1:D:58:ARG:NH1	2.18	0.76
1:A:192:GLN:NE2	2:A:1242:UD2:O7'	2.18	0.76
1:A:49:ASP:OD1	1:A:49:ASP:C	2.28	0.76
1:D:95:ASN:HB2	1:D:98:LEU:HD13	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:PHE:C	1:C:29:MET:HG2	2.11	0.76
1:C:130:ASN:ND2	1:C:149:GLU:O	2.18	0.76
1:A:120:THR:HG22	1:A:208:LEU:HB2	1.66	0.76
1:A:17:TRP:CD1	1:A:47:LEU:CD2	2.68	0.75
1:D:216:GLU:HB3	1:D:227:LEU:HD21	1.69	0.75
1:D:50:GLU:OE1	1:D:58:ARG:CZ	2.33	0.75
1:C:135:TYR:CD2	1:C:145:ILE:HG12	2.20	0.75
1:A:29:MET:CG	1:A:36:ILE:HG13	2.18	0.74
1:C:49:ASP:CG	1:C:49:ASP:O	2.30	0.74
1:A:53:ASN:OD1	1:A:53:ASN:C	2.29	0.74
1:A:53:ASN:O	1:A:56:ILE:HG23	1.87	0.74
1:C:37:GLU:OE2	1:C:57:HIS:NE2	2.20	0.74
1:A:30:GLN:HA	1:A:30:GLN:OE1	1.87	0.74
1:C:48:TYR:CE1	1:C:49:ASP:HB2	2.19	0.74
1:B:118:LEU:HD11	1:B:208:LEU:HG	1.68	0.74
1:C:48:TYR:CD1	1:C:48:TYR:O	2.30	0.74
1:C:133:PHE:HB3	1:C:135:TYR:CE1	2.23	0.74
1:D:158:LEU:HD23	1:D:213:LEU:HD13	1.69	0.74
1:A:50:GLU:OE1	1:A:50:GLU:C	2.32	0.73
1:A:28:PHE:CE2	1:A:29:MET:CE	2.72	0.73
1:D:128:LYS:NZ	1:D:132:GLU:OE1	2.20	0.73
1:C:49:ASP:C	1:C:49:ASP:OD1	2.30	0.73
1:B:28:PHE:CD1	1:B:29:MET:CB	2.72	0.72
1:C:3:ILE:HG22	1:C:90:TYR:HB2	1.71	0.72
1:A:95:ASN:OD1	2:A:1242:UD2:O3B	2.06	0.72
1:C:216:GLU:HB3	1:C:227:LEU:HD21	1.72	0.71
1:B:26:ARG:NH1	1:B:27:TYR:CE1	2.52	0.71
1:B:28:PHE:HD1	1:B:29:MET:CB	2.03	0.71
1:D:36:ILE:O	1:D:55:HIS:ND1	2.24	0.71
1:D:120:THR:HG23	1:D:158:LEU:HD11	1.71	0.71
1:B:28:PHE:CD1	1:B:29:MET:HG2	2.25	0.71
1:D:208:LEU:HB3	1:D:212:TYR:HD2	1.56	0.70
1:D:173:ILE:HG12	1:D:194:LEU:HD11	1.73	0.70
1:D:189:TRP:O	1:D:193:SER:N	2.22	0.70
1:D:92:PHE:HD1	1:D:160:GLY:HA3	1.54	0.70
1:A:139:ASP:OD1	1:A:139:ASP:N	2.24	0.69
1:C:5:ILE:HD11	1:C:28:PHE:CD2	2.27	0.69
1:D:218:TRP:HE3	1:D:220:LEU:HD11	1.57	0.69
1:D:64:LEU:HD13	1:D:69:ASN:HA	1.75	0.69
1:C:145:ILE:HD13	1:C:200:LEU:HD22	1.74	0.69
1:A:54:LYS:CD	1:A:54:LYS:C	2.64	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:ILE:O	2:C:1241:UD2:O2'	2.10	0.69
1:D:40:VAL:O	1:D:59:ILE:HB	1.92	0.69
1:A:26:ARG:HB3	1:A:27:TYR:HD1	1.58	0.69
1:A:28:PHE:HE1	1:A:104:ILE:HG21	1.58	0.68
1:D:29:MET:HG3	1:D:34:PHE:CE1	2.18	0.68
1:D:47:LEU:O	1:D:50:GLU:HB2	1.94	0.68
1:A:188:ILE:CD1	1:A:189:TRP:HD1	1.88	0.68
1:D:50:GLU:CD	1:D:58:ARG:HH12	2.02	0.68
1:A:73:ARG:NH2	2:A:1242:UD2:O6'	2.27	0.68
1:A:198:TYR:O	1:A:202:ASN:ND2	2.27	0.68
1:D:50:GLU:CD	1:D:58:ARG:NH1	2.52	0.68
1:A:84:LEU:HD23	1:A:91:LEU:HD11	1.76	0.67
1:B:78:LEU:HG	1:B:81:LYS:HD2	1.75	0.67
1:A:29:MET:HG3	1:A:36:ILE:HG13	1.77	0.67
1:A:12:LYS:HB3	1:A:240:GLU:HG2	1.77	0.67
1:C:188:ILE:O	1:C:193:SER:OG	2.12	0.67
1:A:54:LYS:HE2	1:A:57:HIS:NE2	2.10	0.67
1:B:46:LYS:C	1:B:47:LEU:HD23	2.20	0.67
1:D:172:THR:O	1:D:175:SER:OG	2.11	0.67
1:D:98:LEU:HD23	1:D:226:ILE:HD11	1.77	0.66
1:B:192:GLN:O	1:B:196:ASN:ND2	2.29	0.66
1:A:53:ASN:CG	1:A:56:ILE:HG22	2.21	0.66
1:C:133:PHE:HB3	1:C:135:TYR:HE1	1.59	0.66
1:C:168:LYS:O	1:C:172:THR:OG1	2.13	0.66
1:A:45:PRO:HG2	1:A:46:LYS:N	2.10	0.66
1:D:70:THR:O	1:D:73:ARG:NE	2.29	0.66
1:A:84:LEU:O	1:A:88:THR:OG1	2.13	0.66
1:B:21:TYR:HE2	1:B:49:ASP:OD2	1.78	0.66
1:C:91:LEU:HG	1:C:163:THR:HG22	1.76	0.66
1:A:50:GLU:OE1	1:A:51:GLU:N	2.29	0.65
1:A:43:ASP:OD1	1:A:44:SER:N	2.29	0.65
1:B:25:GLU:O	1:B:30:GLN:NE2	2.30	0.65
1:C:169:LEU:HA	1:C:198:TYR:HE2	1.62	0.65
1:D:50:GLU:OE1	1:D:58:ARG:NH1	2.29	0.65
1:C:47:LEU:N	1:C:50:GLU:OE2	2.30	0.65
1:A:55:HIS:O	1:A:57:HIS:CD2	2.49	0.65
1:C:192:GLN:OE1	2:C:1241:UD2:O3'	2.12	0.65
1:D:110:PRO:HB2	1:D:116:GLY:HA2	1.78	0.65
1:D:137:ARG:NH1	1:D:147:GLU:OE2	2.30	0.65
1:A:26:ARG:NH1	1:A:27:TYR:OH	2.30	0.64
1:C:70:THR:OG1	1:C:73:ARG:NH2	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:PRO:O	1:B:58:ARG:NH1	2.30	0.64
1:A:53:ASN:OD1	1:A:55:HIS:N	2.29	0.64
1:B:137:ARG:NH1	1:B:146:PRO:O	2.30	0.64
1:C:45:PRO:CD	1:C:46:LYS:N	2.46	0.64
1:C:212:TYR:HA	1:C:226:ILE:HG23	1.80	0.64
1:C:6:LEU:HD11	1:C:91:LEU:HD13	1.79	0.63
1:A:3:ILE:CG1	1:A:29:MET:HE3	2.21	0.63
1:B:134:THR:OG1	1:B:189:TRP:NE1	2.29	0.63
1:A:42:THR:HG23	1:A:43:ASP:N	2.13	0.63
1:B:9:CYS:SG	1:B:14:ASP:HA	2.39	0.63
1:A:214:TYR:HE1	1:A:220:LEU:HB2	1.64	0.63
1:B:84:LEU:O	1:B:88:THR:OG1	2.15	0.63
1:C:213:LEU:HD22	1:C:228:ILE:HD11	1.80	0.63
1:D:122:HIS:NE2	4:D:1238:A2G:O1	2.32	0.63
1:C:31:ASP:OD1	1:C:33:SER:OG	2.15	0.62
1:A:188:ILE:HD11	1:A:193:SER:OG	1.98	0.62
1:B:28:PHE:CD1	1:B:28:PHE:C	2.77	0.62
1:A:29:MET:CB	1:A:36:ILE:CG1	2.77	0.62
1:A:104:ILE:HG22	1:A:105:GLY:H	1.64	0.62
1:A:144:TYR:O	1:A:197:LYS:NZ	2.24	0.62
1:D:2:ARG:HG3	1:D:88:THR:HA	1.82	0.62
1:B:28:PHE:CE1	1:B:29:MET:HG2	2.28	0.62
1:A:69:ASN:HD22	2:A:1242:UD2:C4	2.13	0.62
1:A:188:ILE:HD12	1:A:189:TRP:H	1.59	0.62
1:A:70:THR:OG1	1:A:73:ARG:NH2	2.33	0.62
1:D:40:VAL:CG1	1:D:47:LEU:HD21	2.20	0.62
1:A:3:ILE:HD11	1:A:29:MET:HE3	1.51	0.62
1:A:188:ILE:CD1	1:A:189:TRP:H	2.13	0.61
1:A:188:ILE:CG1	1:A:189:TRP:N	2.62	0.61
1:A:37:GLU:CD	1:A:55:HIS:ND1	2.57	0.61
1:A:45:PRO:CG	1:A:46:LYS:N	2.63	0.61
1:A:22:LEU:HD21	1:A:48:TYR:OH	2.00	0.61
1:C:5:ILE:HD11	1:C:28:PHE:CE2	2.36	0.60
1:C:168:LYS:NZ	1:C:198:TYR:OH	2.34	0.60
1:D:70:THR:O	1:D:73:ARG:CZ	2.49	0.60
1:D:87:GLU:OE1	1:D:87:GLU:N	2.30	0.60
1:A:45:PRO:HG2	1:A:46:LYS:H	1.66	0.60
1:D:17:TRP:HA	1:D:20:PHE:CE1	2.37	0.60
1:A:3:ILE:HG22	1:A:90:TYR:HB2	1.84	0.60
1:D:173:ILE:HD12	1:D:176:TRP:HB2	1.83	0.60
1:D:77:PHE:CZ	1:D:170:CYS:HB2	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:GLU:OE2	1:A:55:HIS:CE1	2.55	0.59
1:A:73:ARG:NH1	1:A:191:ASP:OD2	2.35	0.59
1:A:50:GLU:O	1:A:52:ASN:N	2.30	0.59
1:D:7:TYR:HD2	1:D:40:VAL:HG23	1.65	0.59
1:D:67:PRO:O	1:D:71:LEU:N	2.29	0.59
1:D:7:TYR:CD2	1:D:40:VAL:HG23	2.38	0.59
1:D:115:ASN:OD1	1:D:115:ASN:N	2.35	0.59
1:D:138:ARG:O	1:D:141:SER:OG	2.20	0.59
1:A:73:ARG:HA	1:A:76:ILE:HD12	1.83	0.59
1:B:24:ALA:O	1:B:28:PHE:HB3	2.03	0.59
1:A:50:GLU:C	1:A:51:GLU:CG	2.75	0.59
1:B:177:VAL:O	1:B:180:ASP:HB3	2.01	0.59
1:D:29:MET:HE3	1:D:34:PHE:CE1	2.38	0.58
1:C:106:LYS:HG2	1:C:109:LEU:HD21	1.86	0.58
1:A:192:GLN:NE2	2:A:1242:UD2:H3'	2.16	0.58
1:C:19:ASP:HA	1:C:22:LEU:HD22	1.85	0.58
1:D:211:ALA:HB2	1:D:222:PHE:HB3	1.84	0.58
1:C:216:GLU:OE2	1:C:233:LYS:HE2	2.04	0.58
1:D:77:PHE:HZ	1:D:170:CYS:HB2	1.69	0.58
1:B:147:GLU:OE1	1:B:147:GLU:N	2.36	0.58
1:C:89:ASP:OD1	1:C:89:ASP:N	2.36	0.58
1:D:49:ASP:CG	1:D:53:ASN:HD22	2.12	0.58
1:D:92:PHE:CD1	1:D:160:GLY:HA3	2.37	0.58
1:D:99:LEU:HG	1:D:229:ARG:HG3	1.86	0.58
1:B:136:GLU:OE2	1:B:138:ARG:NH2	2.32	0.58
1:C:232:ASN:HA	1:C:238:GLY:HA2	1.86	0.58
1:D:189:TRP:HB2	1:D:193:SER:HB3	1.84	0.58
1:C:48:TYR:HE1	1:C:49:ASP:HB2	1.60	0.57
1:D:120:THR:HG22	1:D:208:LEU:CB	2.26	0.57
1:C:20:PHE:HE1	1:C:100:PHE:CE1	2.14	0.57
1:B:65:GLY:N	1:B:69:ASN:OD1	2.38	0.57
1:C:81:LYS:HE3	1:C:167:LEU:HD21	1.85	0.57
1:D:19:ASP:HB2	1:D:229:ARG:NH2	2.18	0.57
1:D:186:ILE:HD12	1:D:187:PRO:HD2	1.87	0.57
1:B:145:ILE:HD12	1:B:146:PRO:HD2	1.86	0.56
2:C:1241:UD2:O4'	2:C:1241:UD2:O6'	2.21	0.56
1:A:11:GLY:N	1:A:43:ASP:OD2	2.34	0.56
1:A:55:HIS:O	1:A:57:HIS:NE2	2.38	0.56
1:C:47:LEU:CB	1:C:50:GLU:CD	2.76	0.56
1:C:218:TRP:HE3	1:C:220:LEU:HD11	1.69	0.56
1:A:97:ASN:ND2	1:A:231:LYS:HE3	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:34:PHE:HE2	1:D:36:ILE:HG12	1.70	0.56
1:C:142:THR:HG1	1:C:176:TRP:CG	2.23	0.56
1:D:47:LEU:H	1:D:58:ARG:HH22	1.51	0.56
1:A:137:ARG:NH1	1:A:145:ILE:O	2.33	0.56
1:C:30:GLN:CD	1:C:30:GLN:H	2.14	0.56
1:B:195:ILE:HG23	1:B:199:PHE:CE2	2.40	0.56
1:B:215:PRO:HA	1:B:228:ILE:HB	1.89	0.55
1:A:18:LYS:HG3	1:A:48:TYR:CD2	2.42	0.55
1:B:45:PRO:O	1:B:58:ARG:CZ	2.54	0.55
1:D:29:MET:CB	1:D:34:PHE:CD2	2.69	0.55
1:D:210:PRO:HG2	1:D:222:PHE:CE2	2.42	0.55
1:A:45:PRO:CG	1:A:46:LYS:H	2.20	0.54
1:A:51:GLU:HG3	1:A:52:ASN:OD1	2.07	0.54
1:A:27:TYR:CD1	1:A:27:TYR:N	2.72	0.54
1:A:26:ARG:HB3	1:A:27:TYR:CE1	2.43	0.54
1:D:121:MET:HB2	1:D:210:PRO:HD3	1.90	0.54
1:C:235:GLN:O	1:C:235:GLN:NE2	2.36	0.54
1:A:42:THR:CG2	1:A:43:ASP:N	2.69	0.54
1:C:48:TYR:O	1:C:49:ASP:CB	2.54	0.54
1:D:57:HIS:ND1	1:D:57:HIS:O	2.41	0.54
1:D:166:TYR:O	1:D:169:LEU:HG	2.07	0.53
1:A:10:THR:OG1	1:A:69:ASN:ND2	2.41	0.53
1:A:29:MET:HB2	1:A:36:ILE:CG1	2.38	0.53
1:C:110:PRO:HG2	1:C:161:GLY:HA2	1.90	0.53
1:C:135:TYR:HD2	1:C:145:ILE:HG21	1.73	0.53
1:A:104:ILE:HG22	1:A:105:GLY:N	2.23	0.53
1:B:21:TYR:CE2	1:B:49:ASP:OD2	2.58	0.53
1:B:83:GLN:O	1:B:87:GLU:HG2	2.09	0.53
1:D:47:LEU:HB2	1:D:58:ARG:NH2	2.24	0.53
1:D:130:ASN:HD21	1:D:149:GLU:C	2.16	0.53
1:A:137:ARG:NH1	1:A:146:PRO:O	2.41	0.53
1:C:147:GLU:N	1:C:147:GLU:OE1	2.42	0.53
1:C:48:TYR:HD1	1:C:49:ASP:CB	2.01	0.53
1:A:44:SER:CB	1:A:45:PRO:HD2	2.39	0.53
1:D:9:CYS:SG	1:D:14:ASP:HA	2.48	0.53
1:D:29:MET:CB	1:D:34:PHE:HE2	2.12	0.53
1:D:120:THR:CG2	1:D:158:LEU:HD11	2.38	0.53
1:A:135:TYR:OH	1:A:152:TYR:O	2.22	0.52
1:B:214:TYR:HE2	1:B:224:PRO:HD3	1.74	0.52
1:C:29:MET:SD	1:C:105:GLY:O	2.67	0.52
1:A:6:LEU:HD21	1:A:91:LEU:HD13	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:TYR:HD2	1:A:102:SER:C	2.17	0.52
1:B:37:GLU:OE1	1:B:57:HIS:NE2	2.42	0.52
1:B:135:TYR:OH	1:B:152:TYR:O	2.27	0.52
1:A:17:TRP:NE1	1:A:47:LEU:CD2	2.71	0.52
1:C:199:PHE:HD1	1:C:204:PRO:HD3	1.75	0.52
1:D:43:ASP:OD1	1:D:44:SER:N	2.43	0.52
1:D:188:ILE:HG23	1:D:189:TRP:H	1.74	0.52
1:A:138:ARG:O	1:A:144:TYR:HB2	2.09	0.52
1:B:28:PHE:HD1	1:B:28:PHE:C	2.18	0.52
1:B:210:PRO:HG2	1:B:222:PHE:CZ	2.46	0.51
1:C:133:PHE:HB3	1:C:135:TYR:CD1	2.45	0.51
1:B:210:PRO:HG2	1:B:222:PHE:CE2	2.45	0.51
1:C:48:TYR:O	1:C:49:ASP:HB3	2.09	0.51
1:B:47:LEU:O	1:B:48:TYR:C	2.53	0.51
1:B:68:ASP:O	1:B:72:LYS:HG2	2.10	0.51
1:A:53:ASN:CG	1:A:56:ILE:CG2	2.83	0.51
1:A:191:ASP:OD2	2:A:1242:UD2:O6'	2.21	0.51
1:A:18:LYS:HG3	1:A:48:TYR:CE2	2.45	0.51
1:B:10:THR:OG1	3:B:1239:UDP:O4	2.15	0.51
1:B:125:PHE:HD2	1:B:153:TYR:CD2	2.28	0.51
1:A:169:LEU:HD13	1:A:198:TYR:CD2	2.46	0.51
1:C:7:TYR:CE2	1:C:17:TRP:HE3	2.28	0.51
1:A:29:MET:CB	1:A:36:ILE:CD1	2.74	0.51
1:A:45:PRO:CD	1:A:46:LYS:H	2.24	0.51
1:A:147:GLU:N	1:A:147:GLU:OE1	2.44	0.50
1:B:25:GLU:O	1:B:30:GLN:OE1	2.29	0.50
1:C:48:TYR:O	1:C:49:ASP:OD1	2.30	0.50
1:C:84:LEU:O	1:C:88:THR:OG1	2.14	0.50
1:C:118:LEU:N	1:C:160:GLY:O	2.43	0.50
1:B:126:TYR:OH	1:D:223:GLU:OE2	2.24	0.50
1:B:230:ASP:HB3	1:B:233:LYS:HG3	1.93	0.50
1:D:84:LEU:HD23	1:D:91:LEU:HD21	1.93	0.50
1:A:31:ASP:OD2	1:A:33:SER:OG	2.29	0.50
1:D:172:THR:OG1	1:D:173:ILE:N	2.43	0.50
1:C:195:ILE:HG23	1:C:199:PHE:CE2	2.46	0.50
1:D:1:MET:HG3	1:D:2:ARG:N	2.27	0.50
1:D:214:TYR:HE2	1:D:224:PRO:HD3	1.75	0.50
1:D:135:TYR:OH	1:D:152:TYR:O	2.29	0.50
1:A:2:ARG:HB3	1:A:35:ILE:HD12	1.94	0.50
1:D:146:PRO:HD2	1:D:149:GLU:OE2	2.12	0.50
1:A:51:GLU:HG2	1:A:52:ASN:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ASP:OD1	1:C:43:ASP:N	2.45	0.49
1:A:133:PHE:CD1	1:A:153:TYR:HB2	2.47	0.49
1:B:188:ILE:O	1:B:188:ILE:HD12	2.12	0.49
1:D:40:VAL:HG11	1:D:47:LEU:HD23	1.86	0.49
1:D:214:TYR:HE1	1:D:218:TRP:HB2	1.77	0.49
1:C:48:TYR:CD1	1:C:49:ASP:N	2.80	0.49
1:D:180:ASP:HB3	1:D:185:ILE:O	2.12	0.49
1:A:42:THR:HG22	1:A:44:SER:N	2.25	0.49
1:A:188:ILE:HD12	1:A:188:ILE:C	2.30	0.49
1:A:28:PHE:CE1	1:A:104:ILE:HG21	2.44	0.49
1:A:188:ILE:CG1	1:A:189:TRP:H	2.26	0.49
1:C:65:GLY:N	1:C:69:ASN:OD1	2.46	0.49
1:C:138:ARG:HH12	1:C:140:ALA:HB3	1.76	0.49
1:C:157:GLY:N	2:C:1241:UD2:O6'	2.46	0.49
1:C:215:PRO:HG2	1:C:218:TRP:CD2	2.48	0.49
1:D:97:ASN:O	1:D:229:ARG:N	2.39	0.49
1:D:70:THR:O	1:D:73:ARG:HB3	2.13	0.49
1:D:71:LEU:HA	1:D:73:ARG:NH2	2.28	0.49
1:A:7:TYR:CE1	1:A:96:ALA:HA	2.47	0.49
1:C:95:ASN:H	1:C:98:LEU:HD12	1.78	0.49
1:B:142:THR:HG1	1:B:176:TRP:CG	2.31	0.48
1:A:48:TYR:CD1	1:A:49:ASP:CB	2.92	0.48
1:A:54:LYS:HE2	1:A:57:HIS:CD2	2.48	0.48
1:A:95:ASN:H	1:A:98:LEU:HD12	1.78	0.48
1:D:50:GLU:CD	1:D:58:ARG:NH2	2.67	0.48
1:B:195:ILE:HG23	1:B:199:PHE:HE2	1.77	0.48
1:D:214:TYR:CE2	1:D:224:PRO:HD3	2.48	0.48
1:B:30:GLN:O	1:B:31:ASP:C	2.56	0.48
1:D:35:ILE:HG13	1:D:55:HIS:HE1	1.78	0.48
1:A:38:TYR:C	1:A:39:TYR:HD1	2.21	0.48
1:D:4:GLY:CA	1:D:39:TYR:HE2	2.26	0.48
1:D:50:GLU:CD	1:D:58:ARG:CZ	2.86	0.48
1:D:66:TRP:CZ3	1:D:70:THR:HG21	2.49	0.48
1:D:99:LEU:O	1:D:226:ILE:HD12	2.13	0.48
1:D:79:ARG:HG3	1:D:80:ILE:HG23	1.94	0.48
1:D:194:LEU:C	1:D:194:LEU:HD13	2.39	0.48
1:D:99:LEU:HG	1:D:229:ARG:CG	2.44	0.47
4:D:1238:A2G:O6	4:D:1238:A2G:O4	2.28	0.47
1:D:230:ASP:O	1:D:233:LYS:HG2	2.14	0.47
1:B:23:SER:HB2	1:B:100:PHE:HB2	1.96	0.47
1:D:2:ARG:HA	1:D:35:ILE:HG22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:34:PHE:CE2	1:D:36:ILE:HG12	2.48	0.47
1:C:47:LEU:CA	1:C:50:GLU:OE2	2.62	0.47
1:C:104:ILE:HG13	1:C:105:GLY:H	1.80	0.47
1:D:23:SER:HB2	1:D:100:PHE:HB2	1.96	0.47
1:A:213:LEU:O	1:A:215:PRO:HD3	2.14	0.47
1:D:169:LEU:HB2	1:D:198:TYR:HE2	1.80	0.47
1:D:212:TYR:HA	1:D:226:ILE:HG23	1.97	0.47
1:B:78:LEU:O	1:B:81:LYS:HD3	2.14	0.47
1:D:8:ILE:HG23	1:D:93:PHE:HE2	1.80	0.47
1:D:11:GLY:H	1:D:63:ASN:HD21	1.63	0.47
1:A:53:ASN:OD1	1:A:53:ASN:O	2.32	0.47
1:A:145:ILE:HG12	1:A:197:LYS:HE2	1.97	0.47
1:A:151:ARG:NH2	1:A:152:TYR:OH	2.48	0.47
1:B:191:ASP:N	1:B:191:ASP:OD1	2.47	0.47
1:C:18:LYS:HG3	1:C:48:TYR:CD2	2.49	0.47
1:C:138:ARG:NH1	1:C:185:ILE:HD12	2.30	0.47
1:C:19:ASP:O	1:C:23:SER:OG	2.31	0.47
1:D:7:TYR:CE1	1:D:96:ALA:HA	2.50	0.47
1:D:101:THR:OG1	1:D:225:ILE:O	2.29	0.47
1:B:231:LYS:NZ	3:B:1239:UDP:O1B	2.48	0.47
1:B:74:PHE:O	1:B:78:LEU:HD13	2.15	0.47
1:B:90:TYR:HE2	1:B:109:LEU:HD13	1.80	0.46
1:C:208:LEU:HD23	1:C:212:TYR:CD2	2.50	0.46
1:A:59:ILE:HD13	1:A:80:ILE:HD13	1.98	0.46
1:D:117:LEU:HD13	1:D:204:PRO:HG3	1.98	0.46
1:B:73:ARG:HA	1:B:76:ILE:HD12	1.98	0.46
1:D:208:LEU:HB3	1:D:212:TYR:CD2	2.42	0.46
1:D:3:ILE:HD12	1:D:34:PHE:HE1	1.80	0.46
1:B:1:MET:HB3	1:B:89:ASP:OD2	2.15	0.46
1:B:187:PRO:HG2	1:B:194:LEU:HD11	1.97	0.46
1:D:134:THR:OG1	1:D:153:TYR:OH	2.25	0.46
1:A:34:PHE:CZ	1:A:106:LYS:HG2	2.50	0.46
1:A:47:LEU:O	1:A:50:GLU:HB2	2.15	0.46
1:A:173:ILE:O	1:A:177:VAL:HG12	2.16	0.46
1:C:2:ARG:NH2	1:C:86:ARG:O	2.41	0.46
1:C:28:PHE:O	1:C:28:PHE:CD1	2.68	0.46
1:C:138:ARG:CZ	1:C:141:SER:HB2	2.45	0.46
1:C:145:ILE:HD12	1:C:146:PRO:HD2	1.97	0.46
1:D:169:LEU:HB2	1:D:198:TYR:CE2	2.51	0.46
1:B:199:PHE:CD1	1:B:204:PRO:HG3	2.50	0.46
1:C:7:TYR:HD1	1:C:8:ILE:N	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:LEU:HD23	1:C:91:LEU:HA	1.81	0.46
1:A:30:GLN:OE1	1:A:30:GLN:CA	2.59	0.45
1:A:98:LEU:HD22	1:A:226:ILE:HG22	1.96	0.45
1:B:176:TRP:O	1:B:180:ASP:HB2	2.16	0.45
1:D:173:ILE:HD13	1:D:176:TRP:CE3	2.51	0.45
1:D:220:LEU:HB3	1:D:222:PHE:CD2	2.52	0.45
1:A:235:GLN:H	1:A:235:GLN:HG2	1.53	0.45
1:C:8:ILE:HG22	1:C:93:PHE:HE1	1.82	0.45
1:D:64:LEU:HD12	1:D:64:LEU:H	1.81	0.45
1:B:25:GLU:O	1:B:30:GLN:CD	2.60	0.45
1:B:136:GLU:HB3	1:B:143:ALA:O	2.17	0.45
1:A:54:LYS:HG3	1:A:55:HIS:N	2.29	0.45
1:C:20:PHE:CE1	1:C:100:PHE:HE1	2.15	0.45
1:C:40:VAL:HG21	1:C:47:LEU:HD11	1.99	0.45
1:B:8:ILE:HD11	1:B:77:PHE:CE1	2.52	0.45
1:A:2:ARG:HB3	1:A:35:ILE:CD1	2.47	0.45
1:C:27:TYR:CE1	1:C:103:PRO:HD3	2.52	0.45
1:D:134:THR:HG1	1:D:153:TYR:HH	1.57	0.45
1:B:7:TYR:CE2	1:B:96:ALA:HA	2.52	0.45
1:B:222:PHE:O	1:D:221:PRO:HB3	2.16	0.45
1:C:73:ARG:HB2	1:C:77:PHE:CZ	2.51	0.45
1:D:20:PHE:HB3	1:D:99:LEU:HD23	1.99	0.45
1:D:76:ILE:HD12	1:D:76:ILE:HA	1.75	0.45
1:A:17:TRP:NE1	1:A:47:LEU:HD22	2.32	0.44
1:A:236:TYR:C	1:A:237:GLY:O	2.57	0.44
1:B:18:LYS:HE3	1:B:48:TYR:CE1	2.52	0.44
1:B:43:ASP:OD1	1:B:44:SER:N	2.51	0.44
1:B:203:PRO:HA	1:B:204:PRO:HD3	1.93	0.44
1:D:4:GLY:HA3	1:D:39:TYR:HE2	1.82	0.44
1:B:90:TYR:HD2	1:B:109:LEU:HD22	1.83	0.44
1:C:28:PHE:HD1	1:C:104:ILE:HD11	1.82	0.44
1:C:49:ASP:OD1	1:C:56:ILE:CD1	2.66	0.44
1:A:55:HIS:O	1:A:57:HIS:CE1	2.70	0.44
1:B:135:TYR:CD2	1:B:145:ILE:HG12	2.53	0.44
1:D:152:TYR:HE2	1:D:207:THR:HG21	1.82	0.44
1:A:17:TRP:CG	1:A:47:LEU:CD2	3.00	0.44
1:C:130:ASN:HA	1:C:133:PHE:CD2	2.53	0.44
1:D:16:PHE:HE1	1:D:236:TYR:CD2	2.36	0.44
1:A:3:ILE:HD11	1:A:29:MET:HG3	1.98	0.44
1:A:179:ARG:O	1:A:182:THR:HG22	2.17	0.44
1:C:9:CYS:SG	1:C:14:ASP:HA	2.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LEU:HD23	1:C:91:LEU:HD11	2.00	0.44
1:C:138:ARG:CZ	1:C:185:ILE:HD12	2.48	0.44
1:A:214:TYR:CE1	1:A:220:LEU:HB2	2.50	0.44
1:B:26:ARG:NH1	1:B:27:TYR:OH	2.50	0.43
1:B:214:TYR:O	1:B:228:ILE:N	2.43	0.43
1:D:94:PHE:HD2	1:D:98:LEU:HD23	1.82	0.43
1:D:173:ILE:CG1	1:D:194:LEU:HD11	2.45	0.43
1:A:62:LYS:HD2	1:A:62:LYS:HA	1.68	0.43
1:A:208:LEU:HB3	1:A:212:TYR:CD2	2.54	0.43
1:D:7:TYR:CE1	1:D:20:PHE:HZ	2.36	0.43
1:D:24:ALA:HA	1:D:100:PHE:CE2	2.53	0.43
1:A:34:PHE:HZ	1:A:106:LYS:HG2	1.82	0.43
1:D:91:LEU:HD13	1:D:163:THR:OG1	2.18	0.43
1:B:180:ASP:OD1	1:B:185:ILE:HG13	2.19	0.43
1:C:34:PHE:CD1	1:C:34:PHE:C	2.96	0.43
1:C:173:ILE:HD12	1:C:194:LEU:HD23	2.01	0.43
1:D:76:ILE:O	1:D:79:ARG:HG2	2.19	0.43
1:A:168:LYS:HD2	1:A:168:LYS:HA	1.67	0.43
1:B:214:TYR:CE2	1:B:224:PRO:HD3	2.53	0.43
1:A:17:TRP:CG	1:A:47:LEU:HD21	2.53	0.43
1:A:91:LEU:HD23	1:A:91:LEU:HA	1.84	0.43
1:A:48:TYR:CD1	1:A:49:ASP:CA	3.02	0.43
1:A:59:ILE:CD1	1:A:80:ILE:HD13	2.48	0.43
1:C:17:TRP:CD1	1:C:48:TYR:HB3	2.54	0.43
1:C:142:THR:HG21	1:C:176:TRP:O	2.19	0.43
1:A:17:TRP:CE2	1:A:47:LEU:HD22	2.53	0.42
1:C:133:PHE:CB	1:C:135:TYR:HE1	2.28	0.42
1:D:25:GLU:OE2	1:D:38:TYR:OH	2.25	0.42
1:D:169:LEU:O	1:D:173:ILE:HG22	2.19	0.42
1:A:115:ASN:ND2	1:A:203:PRO:O	2.52	0.42
1:C:16:PHE:HB3	1:C:96:ALA:O	2.19	0.42
1:C:27:TYR:O	1:C:28:PHE:C	2.58	0.42
1:D:20:PHE:O	1:D:24:ALA:N	2.43	0.42
1:D:92:PHE:CE1	1:D:118:LEU:HD21	2.54	0.42
1:A:48:TYR:HA	1:A:49:ASP:HA	1.73	0.42
1:B:28:PHE:HD2	1:B:100:PHE:CZ	2.37	0.42
1:D:17:TRP:HB2	1:D:21:TYR:CE2	2.54	0.42
1:A:40:VAL:CG1	1:A:47:LEU:CD1	2.82	0.42
1:D:7:TYR:HE1	1:D:96:ALA:HA	1.84	0.42
1:D:173:ILE:HD12	1:D:173:ILE:O	2.19	0.42
1:B:180:ASP:OD2	1:B:185:ILE:O	2.38	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:TYR:HE2	1:C:20:PHE:CD2	2.37	0.42
1:B:21:TYR:CZ	1:B:25:GLU:OE1	2.73	0.42
1:C:2:ARG:HH22	1:C:86:ARG:C	2.24	0.42
1:C:173:ILE:O	1:C:176:TRP:HB2	2.20	0.42
1:D:32:GLN:C	1:D:34:PHE:H	2.28	0.42
1:D:153:TYR:CZ	1:D:192:GLN:NE2	2.88	0.42
1:D:208:LEU:HD22	1:D:212:TYR:CE2	2.54	0.42
1:C:45:PRO:HD2	1:C:46:LYS:N	2.22	0.42
1:B:107:GLU:HG2	1:B:108:ILE:HG23	2.02	0.42
1:B:135:TYR:O	1:B:188:ILE:HD11	2.20	0.42
1:B:152:TYR:CD2	1:B:207:THR:HG21	2.54	0.42
1:C:2:ARG:HA	1:C:35:ILE:O	2.19	0.42
1:C:95:ASN:N	1:C:98:LEU:HD12	2.34	0.42
1:C:225:ILE:HG23	1:C:226:ILE:H	1.85	0.42
1:D:4:GLY:O	1:D:91:LEU:HA	2.20	0.42
1:A:104:ILE:CG2	1:A:105:GLY:H	2.33	0.42
1:A:212:TYR:O	1:A:226:ILE:HB	2.19	0.42
1:C:27:TYR:HE1	1:C:103:PRO:HD3	1.84	0.42
1:D:48:TYR:O	1:D:49:ASP:CB	2.67	0.42
1:C:206:ILE:HG13	1:C:208:LEU:HD12	2.02	0.41
1:A:143:ALA:HA	1:A:197:LYS:HE3	2.02	0.41
1:C:195:ILE:HG23	1:C:199:PHE:CD2	2.55	0.41
1:B:18:LYS:HE3	1:B:48:TYR:CZ	2.56	0.41
1:D:71:LEU:CD2	1:D:186:ILE:HD11	2.50	0.41
1:C:1:MET:HE3	1:C:1:MET:HB2	1.95	0.41
1:D:39:TYR:HD1	1:D:57:HIS:CE1	2.38	0.41
1:D:158:LEU:HD12	1:D:158:LEU:C	2.46	0.41
1:A:107:GLU:HG2	1:A:108:ILE:HG23	2.02	0.41
1:A:126:TYR:CE2	1:B:219:LEU:HD21	2.56	0.41
1:A:202:ASN:HA	1:A:203:PRO:HD2	1.93	0.41
1:A:222:PHE:HA	1:B:216:GLU:HG3	2.02	0.41
1:B:90:TYR:CD2	1:B:109:LEU:HD22	2.56	0.41
1:B:170:CYS:O	1:B:174:CYS:HB3	2.21	0.41
1:C:121:MET:HE3	1:C:152:TYR:CD1	2.54	0.41
1:B:235:GLN:H	1:B:235:GLN:HG2	1.73	0.41
2:C:1241:UD2:H6	2:C:1241:UD2:H2B	1.50	0.41
1:A:130:ASN:HD21	1:A:149:GLU:C	2.28	0.41
1:B:89:ASP:C	1:B:163:THR:HG1	2.24	0.41
1:B:110:PRO:HG2	1:B:161:GLY:HA2	2.03	0.41
1:C:194:LEU:HD22	1:C:194:LEU:H	1.85	0.41
1:A:136:GLU:HB3	1:A:143:ALA:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:LYS:HE2	1:B:72:LYS:HB3	1.85	0.41
1:B:90:TYR:CE2	1:B:109:LEU:HD13	2.55	0.41
1:B:142:THR:HG21	1:B:176:TRP:HB3	2.01	0.41
1:C:135:TYR:HB3	1:C:145:ILE:HG23	2.03	0.41
1:D:7:TYR:CE1	1:D:20:PHE:CZ	3.08	0.41
1:D:39:TYR:CD1	1:D:57:HIS:NE2	2.89	0.41
1:D:173:ILE:CD1	1:D:194:LEU:HD21	2.51	0.41
1:D:190:HIS:O	1:D:194:LEU:HB3	2.21	0.41
1:A:32:GLN:O	1:A:33:SER:C	2.61	0.41
1:B:94:PHE:HD2	1:B:98:LEU:HD12	1.86	0.41
1:B:120:THR:HG21	1:B:213:LEU:HD13	2.03	0.40
1:B:134:THR:HG1	1:B:189:TRP:CD1	2.39	0.40
1:A:64:LEU:HB2	1:A:69:ASN:OD1	2.21	0.40
1:B:158:LEU:CB	1:B:213:LEU:HD11	2.51	0.40
1:C:50:GLU:H	1:C:50:GLU:HG2	1.40	0.40
1:D:29:MET:O	1:D:34:PHE:CD2	2.74	0.40
1:D:198:TYR:HD1	1:D:198:TYR:O	2.04	0.40
1:D:66:TRP:CG	1:D:67:PRO:HA	2.57	0.40
1:B:168:LYS:HA	1:B:171:THR:HG22	2.04	0.40
1:C:47:LEU:CD2	1:C:56:ILE:CG2	2.87	0.40
1:C:115:ASN:OD1	1:C:116:GLY:N	2.54	0.40
1:A:95:ASN:N	1:A:98:LEU:HD12	2.35	0.40
1:D:30:GLN:O	1:D:31:ASP:C	2.63	0.40
1:D:102:SER:HA	1:D:103:PRO:HD3	1.92	0.40
1:D:144:TYR:CE2	1:D:146:PRO:HG3	2.56	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	239/246 (97%)	222 (93%)	17 (7%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	236/246 (96%)	219 (93%)	16 (7%)	1 (0%)	30	59
1	C	238/246 (97%)	219 (92%)	19 (8%)	0	100	100
1	D	234/246 (95%)	218 (93%)	15 (6%)	1 (0%)	30	59
All	All	947/984 (96%)	878 (93%)	67 (7%)	2 (0%)	43	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	146	PRO
1	B	146	PRO

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	216/221 (98%)	182 (84%)	34 (16%)	2	10
1	B	213/221 (96%)	192 (90%)	21 (10%)	7	26
1	C	215/221 (97%)	176 (82%)	39 (18%)	2	7
1	D	213/221 (96%)	193 (91%)	20 (9%)	8	28
All	All	857/884 (97%)	743 (87%)	114 (13%)	4	15

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	ILE
1	A	6	LEU
1	A	28	PHE
1	A	29	MET
1	A	30	GLN
1	A	32	GLN
1	A	33	SER
1	A	35	ILE

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Mol	Chain	Res	Type
1	A	39	TYR
1	A	42	THR
1	A	44	SER
1	A	47	LEU
1	A	50	GLU
1	A	51	GLU
1	A	53	ASN
1	A	54	LYS
1	A	56	ILE
1	A	66	TRP
1	A	102	SER
1	A	121	MET
1	A	139	ASP
1	A	158	LEU
1	A	175	SER
1	A	177	VAL
1	A	186	ILE
1	A	187	PRO
1	A	190	HIS
1	A	195	ILE
1	A	202	ASN
1	A	209	SER
1	A	235	GLN
1	A	239	HIS
1	A	241	LEU
1	B	10	THR
1	B	15	ILE
1	B	18	LYS
1	B	25	GLU
1	B	40	VAL
1	B	47	LEU
1	B	51	GLU
1	B	56	ILE
1	B	81	LYS
1	B	89	ASP
1	B	98	LEU
1	B	134	THR
1	B	142	THR
1	B	145	ILE
1	B	163	THR
1	B	174	CYS
1	B	177	VAL

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Mol	Chain	Res	Type
1	B	182	THR
1	B	189	TRP
1	B	195	ILE
1	B	206	ILE
1	C	1	MET
1	C	8	ILE
1	C	14	ASP
1	C	22	LEU
1	C	29	MET
1	C	35	ILE
1	C	40	VAL
1	C	43	ASP
1	C	44	SER
1	C	45	PRO
1	C	46	LYS
1	C	47	LEU
1	C	48	TYR
1	C	49	ASP
1	C	50	GLU
1	C	64	LEU
1	C	89	ASP
1	C	104	ILE
1	C	106	LYS
1	C	109	LEU
1	C	138	ARG
1	C	142	THR
1	C	145	ILE
1	C	149	GLU
1	C	158	LEU
1	C	163	THR
1	C	164	LYS
1	C	171	THR
1	C	172	THR
1	C	173	ILE
1	C	183	ASN
1	C	185	ILE
1	C	190	HIS
1	C	194	LEU
1	C	202	ASN
1	C	223	GLU
1	C	225	ILE
1	C	226	ILE

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Mol	Chain	Res	Type
1	C	235	GLN
1	D	8	ILE
1	D	21	TYR
1	D	29	MET
1	D	30	GLN
1	D	40	VAL
1	D	47	LEU
1	D	56	ILE
1	D	62	LYS
1	D	64	LEU
1	D	76	ILE
1	D	117	LEU
1	D	118	LEU
1	D	145	ILE
1	D	162	CYS
1	D	174	CYS
1	D	177	VAL
1	D	192	GLN
1	D	194	LEU
1	D	202	ASN
1	D	226	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	ASN
1	B	61	GLN
1	B	235	GLN
1	C	202	ASN
1	D	53	ASN
1	D	69	ASN
1	D	192	GLN
1	D	196	ASN
1	D	202	ASN

5.3.3 RNA ⓘ

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](#)

6 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	$\# Z > 2$
3	UDP	D	1237	-	25,26,26	1.19	3 (12%)	38,40,40	1.75	7 (18%)
4	A2G	D	1238	-	15,15,15	0.43	0	21,21,21	0.38	0
2	UD2	A	1242	-	40,41,41	1.04	5 (12%)	59,62,62	1.58	11 (18%)
3	UDP	B	1239	-	25,26,26	1.34	4 (16%)	38,40,40	1.85	8 (21%)
4	A2G	B	1240	-	15,15,15	0.42	0	21,21,21	0.65	0
2	UD2	C	1241	-	40,41,41	1.01	4 (10%)	59,62,62	1.62	11 (18%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UDP	D	1237	-	-	2/16/32/32	0/2/2/2
4	A2G	D	1238	-	-	4/6/26/26	0/1/1/1
2	UD2	A	1242	-	-	16/26/63/63	0/3/3/3
3	UDP	B	1239	-	-	11/16/32/32	0/2/2/2
4	A2G	B	1240	-	-	3/6/26/26	0/1/1/1
2	UD2	C	1241	-	-	10/26/63/63	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1239	UDP	C4-N3	-3.24	1.33	1.38
3	B	1239	UDP	C2-N1	3.02	1.43	1.38
2	A	1242	UD2	C4-N3	-2.71	1.34	1.38
2	C	1241	UD2	C4-N3	-2.69	1.34	1.38
3	D	1237	UDP	C2-N1	2.65	1.42	1.38
3	B	1239	UDP	C2-N3	-2.58	1.33	1.38
2	A	1242	UD2	C2-N1	2.56	1.42	1.38
3	D	1237	UDP	C4-N3	-2.36	1.34	1.38
3	D	1237	UDP	C2-N3	-2.36	1.33	1.38
2	C	1241	UD2	C2-N1	2.29	1.42	1.38
2	C	1241	UD2	C5-C4	-2.13	1.39	1.43
2	A	1242	UD2	C5-C4	-2.12	1.39	1.43
2	A	1242	UD2	C2-N3	-2.09	1.34	1.38
2	C	1241	UD2	C2-N3	-2.08	1.34	1.38
2	A	1242	UD2	PA-O3A	2.07	1.61	1.59
3	B	1239	UDP	C5-C4	-2.03	1.39	1.43

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1237	UDP	C4-N3-C2	-5.11	120.27	126.61
3	B	1239	UDP	C1'-N1-C2	4.63	125.91	117.59
3	B	1239	UDP	C4-N3-C2	-4.60	120.91	126.61
3	B	1239	UDP	N3-C2-N1	4.54	120.80	114.89
2	C	1241	UD2	C4-N3-C2	-4.46	121.07	126.61
2	A	1242	UD2	C4-N3-C2	-4.40	121.15	126.61
3	D	1237	UDP	C5-C4-N3	4.35	120.89	114.80
3	D	1237	UDP	N3-C2-N1	4.25	120.42	114.89
2	C	1241	UD2	N3-C2-N1	4.13	120.27	114.89
2	A	1242	UD2	C1B-N1-C2	4.07	124.90	117.59
2	C	1241	UD2	O5'-C5'-C4'	4.05	117.00	109.70
2	A	1242	UD2	N3-C2-N1	3.84	119.89	114.89
2	A	1242	UD2	C5-C4-N3	3.78	120.09	114.80
3	B	1239	UDP	C5-C4-N3	3.66	119.92	114.80
3	B	1239	UDP	O2-C2-N3	-3.64	114.77	121.49
2	C	1241	UD2	C5-C4-N3	3.58	119.81	114.80
2	A	1242	UD2	C2B-C1B-N1	-3.36	103.89	113.25
3	D	1237	UDP	C3'-C2'-C1'	3.28	107.68	101.46
3	D	1237	UDP	O4-C4-C5	-3.14	119.75	125.16
2	C	1241	UD2	C3'-C4'-C5'	3.08	115.82	110.23
2	A	1242	UD2	O4-C4-C5	-3.07	119.88	125.16
2	C	1241	UD2	O4-C4-C5	-2.99	120.00	125.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1241	UD2	C2'-N2'-C7'	-2.93	116.24	123.11
2	A	1242	UD2	C3'-C2'-N2'	2.65	115.50	110.62
2	C	1241	UD2	C3B-C2B-C1B	2.62	106.42	101.46
3	D	1237	UDP	O2-C2-N1	-2.47	119.58	122.80
2	C	1241	UD2	C3'-C2'-N2'	2.40	115.04	110.62
3	B	1239	UDP	C1'-N1-C6	-2.40	115.65	120.78
2	C	1241	UD2	C1'-O5'-C5'	2.38	118.37	113.72
2	A	1242	UD2	C1B-N1-C6	-2.27	115.92	120.78
3	B	1239	UDP	O4-C4-C5	-2.26	121.26	125.16
2	C	1241	UD2	C1B-N1-C2	2.21	121.56	117.59
2	A	1242	UD2	C4'-C3'-C2'	2.16	113.54	110.40
2	A	1242	UD2	O4B-C1B-N1	2.12	113.17	108.36
3	B	1239	UDP	C6-N1-C2	-2.12	118.42	121.00
2	A	1242	UD2	O2-C2-N3	-2.09	117.64	121.49
3	D	1237	UDP	C1'-N1-C2	2.02	121.22	117.59

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1242	UD2	C3'-C2'-N2'-C7'
2	A	1242	UD2	C8'-C7'-N2'-C2'
2	A	1242	UD2	O7'-C7'-N2'-C2'
2	A	1242	UD2	C1'-O1'-PB-O1B
2	A	1242	UD2	O4B-C1B-N1-C2
2	A	1242	UD2	O4B-C1B-N1-C6
2	A	1242	UD2	C5B-O5B-PA-O1A
2	A	1242	UD2	C5B-O5B-PA-O2A
2	A	1242	UD2	C5B-O5B-PA-O3A
2	C	1241	UD2	C1'-O1'-PB-O3A
2	C	1241	UD2	O4B-C1B-N1-C2
2	C	1241	UD2	O4B-C1B-N1-C6
2	C	1241	UD2	C5B-O5B-PA-O2A
3	B	1239	UDP	C3'-C4'-C5'-O5'
3	B	1239	UDP	O4'-C4'-C5'-O5'
3	B	1239	UDP	C5'-O5'-PA-O2A
3	B	1239	UDP	C5'-O5'-PA-O3A
4	B	1240	A2G	O7-C7-N2-C2
4	B	1240	A2G	C8-C7-N2-C2
4	D	1238	A2G	O7-C7-N2-C2
4	D	1238	A2G	C8-C7-N2-C2
2	C	1241	UD2	C2B-C1B-N1-C6

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Mol	Chain	Res	Type	Atoms
2	C	1241	UD2	C3B-C4B-C5B-O5B
2	C	1241	UD2	O4B-C4B-C5B-O5B
2	A	1242	UD2	O4B-C4B-C5B-O5B
2	A	1242	UD2	C4'-C5'-C6'-O6'
4	B	1240	A2G	O5-C5-C6-O6
2	C	1241	UD2	C2B-C1B-N1-C2
2	C	1241	UD2	O5'-C5'-C6'-O6'
3	B	1239	UDP	C2'-C1'-N1-C2
2	A	1242	UD2	O5'-C5'-C6'-O6'
3	B	1239	UDP	C4'-C5'-O5'-PA
3	B	1239	UDP	C2'-C1'-N1-C6
4	D	1238	A2G	C4-C5-C6-O6
2	A	1242	UD2	C3B-C4B-C5B-O5B
2	A	1242	UD2	C1'-O1'-PB-O2B
2	C	1241	UD2	C1'-O1'-PB-O2B
3	D	1237	UDP	PB-O3A-PA-O1A
3	B	1239	UDP	O4'-C1'-N1-C2
4	D	1238	A2G	O5-C5-C6-O6
3	B	1239	UDP	C5'-O5'-PA-O1A
3	B	1239	UDP	O4'-C1'-N1-C6
2	A	1242	UD2	O5'-C1'-O1'-PB
3	D	1237	UDP	PB-O3A-PA-O2A
2	A	1242	UD2	PB-O3A-PA-O1A
3	B	1239	UDP	PB-O3A-PA-O2A

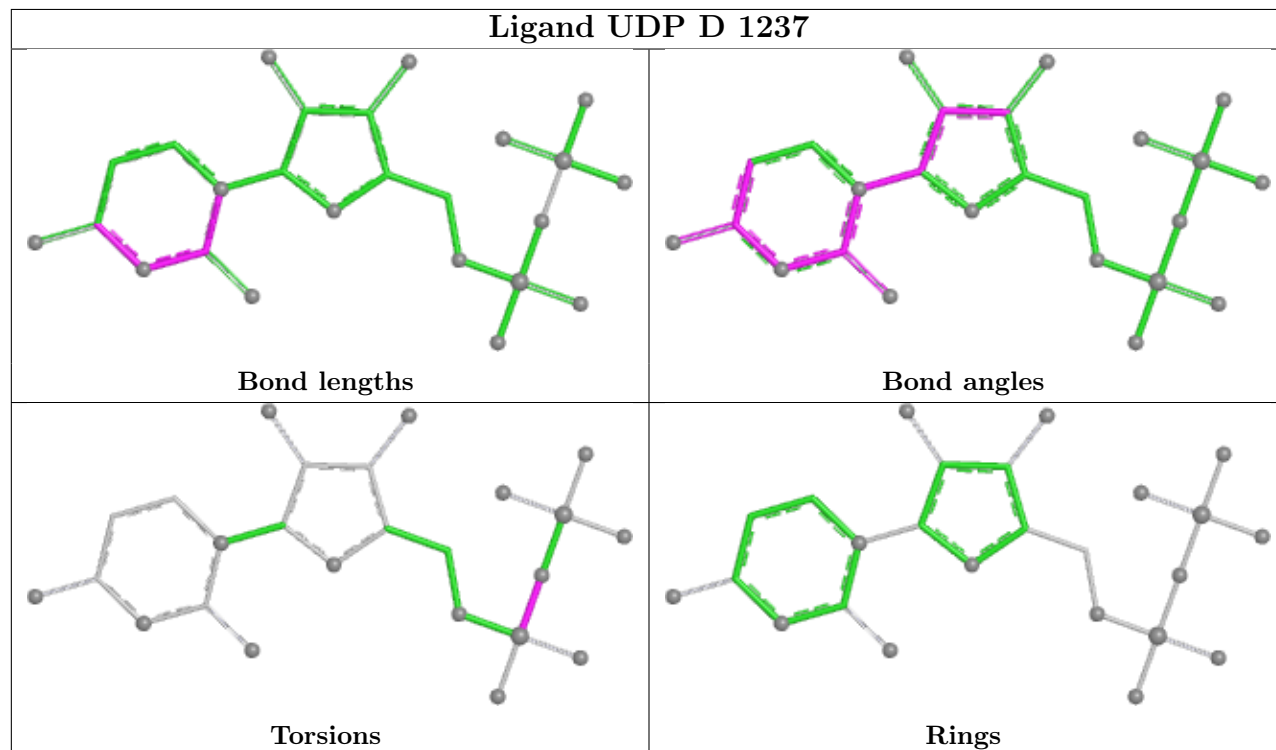
There are no ring outliers.

5 monomers are involved in 20 short contacts:

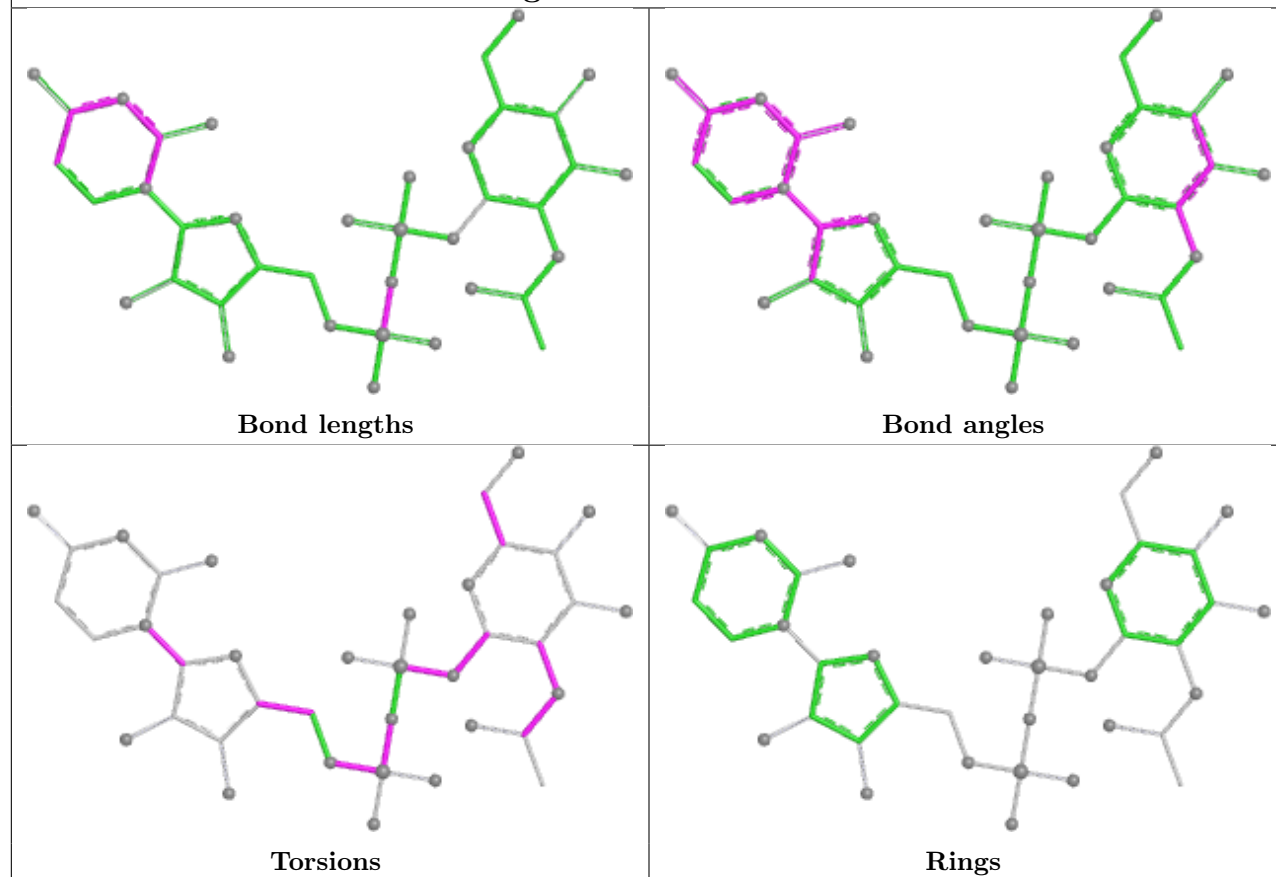
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1237	UDP	2	0
4	D	1238	A2G	2	0
2	A	1242	UD2	9	0
3	B	1239	UDP	2	0
2	C	1241	UD2	5	0

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

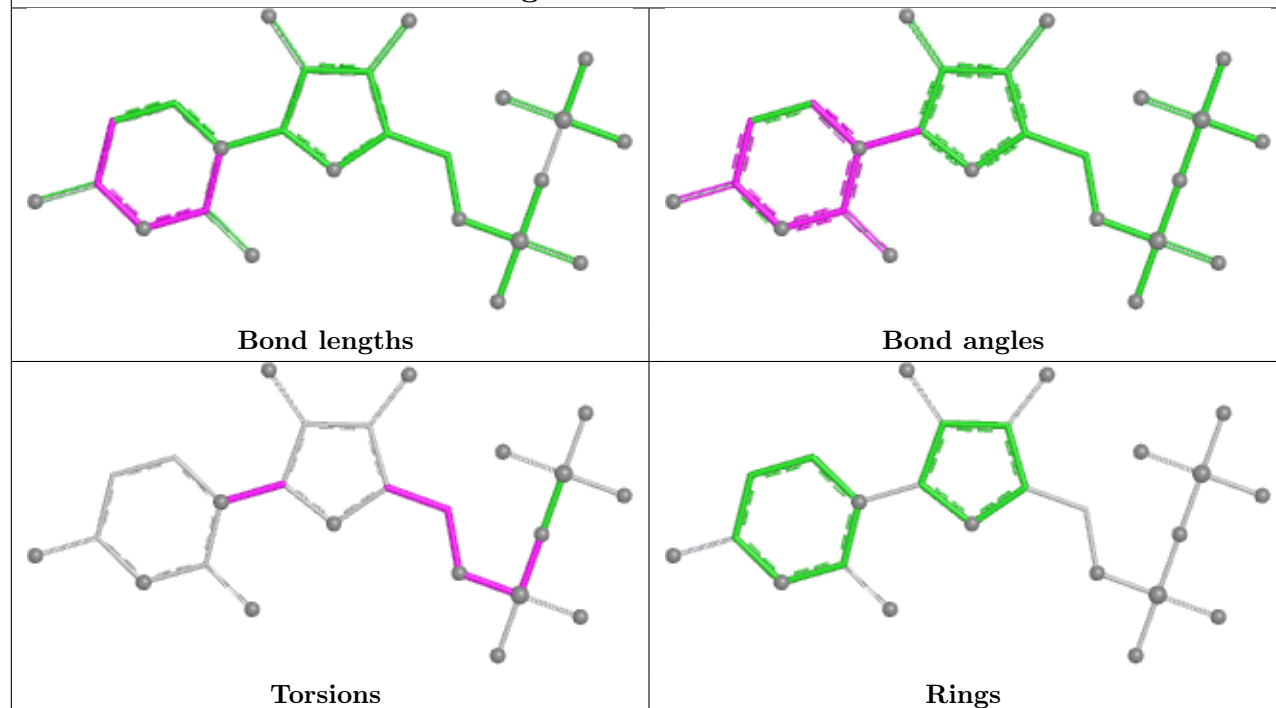
in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

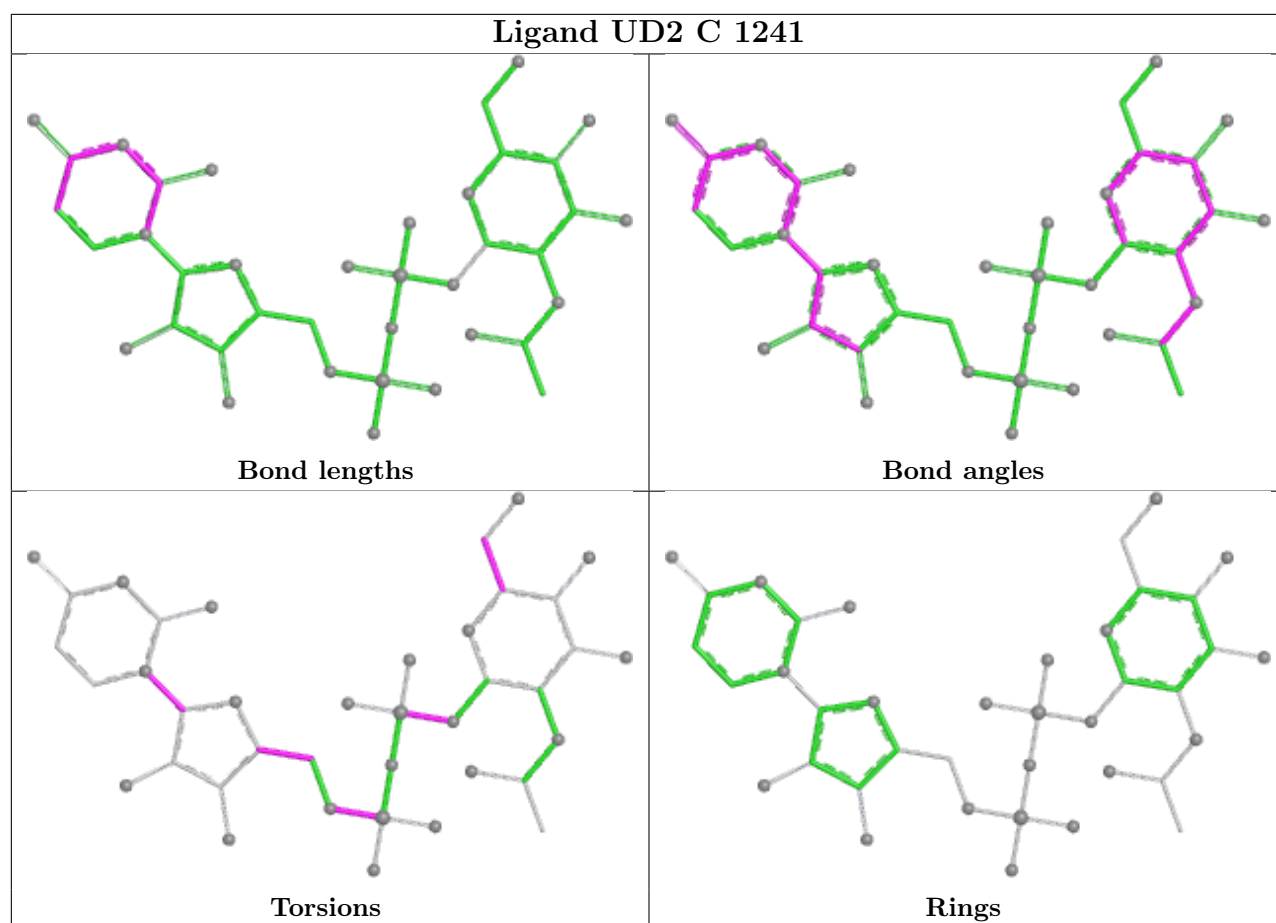


Ligand UD2 A 1242



Ligand UDP B 1239





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	241/246 (97%)	0.08	2 (0%) 82 71	23, 51, 79, 94	2 (0%)
1	B	238/246 (96%)	0.17	5 (2%) 63 48	26, 63, 84, 95	3 (1%)
1	C	240/246 (97%)	0.38	7 (2%) 53 39	25, 96, 148, 184	3 (1%)
1	D	236/246 (95%)	1.09	39 (16%) 4 6	46, 117, 199, 358	3 (1%)
All	All	955/984 (97%)	0.43	53 (5%) 30 23	23, 74, 160, 358	11 (1%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	108	ILE	8.2
1	D	91	LEU	5.6
1	D	117	LEU	5.3
1	D	166	TYR	4.5
1	A	29	MET	4.4
1	D	5	ILE	4.0
1	D	39	TYR	4.0
1	D	29	MET	3.8
1	D	163	THR	3.6
1	B	123	PRO	3.5
1	D	118	LEU	3.5
1	D	85	GLU	3.4
1	D	84	LEU	3.3
1	D	4	GLY	3.2
1	D	119	GLY	3.2
1	D	28	PHE	3.2
1	D	3	ILE	3.1
1	D	36	ILE	3.1
1	D	6	LEU	3.0
1	B	200	LEU	2.9
1	D	167	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	76	ILE	2.8
1	D	104	ILE	2.6
1	D	111	PRO	2.6
1	D	160	GLY	2.6
1	D	116	GLY	2.5
1	C	71	LEU	2.5
1	D	8	ILE	2.4
1	C	47	LEU	2.3
1	D	90	TYR	2.3
1	D	115	ASN	2.3
1	D	17	TRP	2.3
1	D	38	TYR	2.3
1	C	159	SER	2.3
1	C	191	ASP	2.3
1	D	34	PHE	2.3
1	D	206	ILE	2.2
1	A	148	GLY	2.2
1	D	114	SER	2.2
1	D	159	SER	2.2
1	C	178	ASP	2.2
1	B	124	GLY	2.2
1	D	168	LYS	2.2
1	D	171	THR	2.2
1	D	92	PHE	2.2
1	D	56	ILE	2.1
1	C	240	GLU	2.1
1	D	158	LEU	2.1
1	D	222	PHE	2.1
1	D	88	THR	2.1
1	D	161	GLY	2.1
1	B	186	ILE	2.0
1	B	81	LYS	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](#)

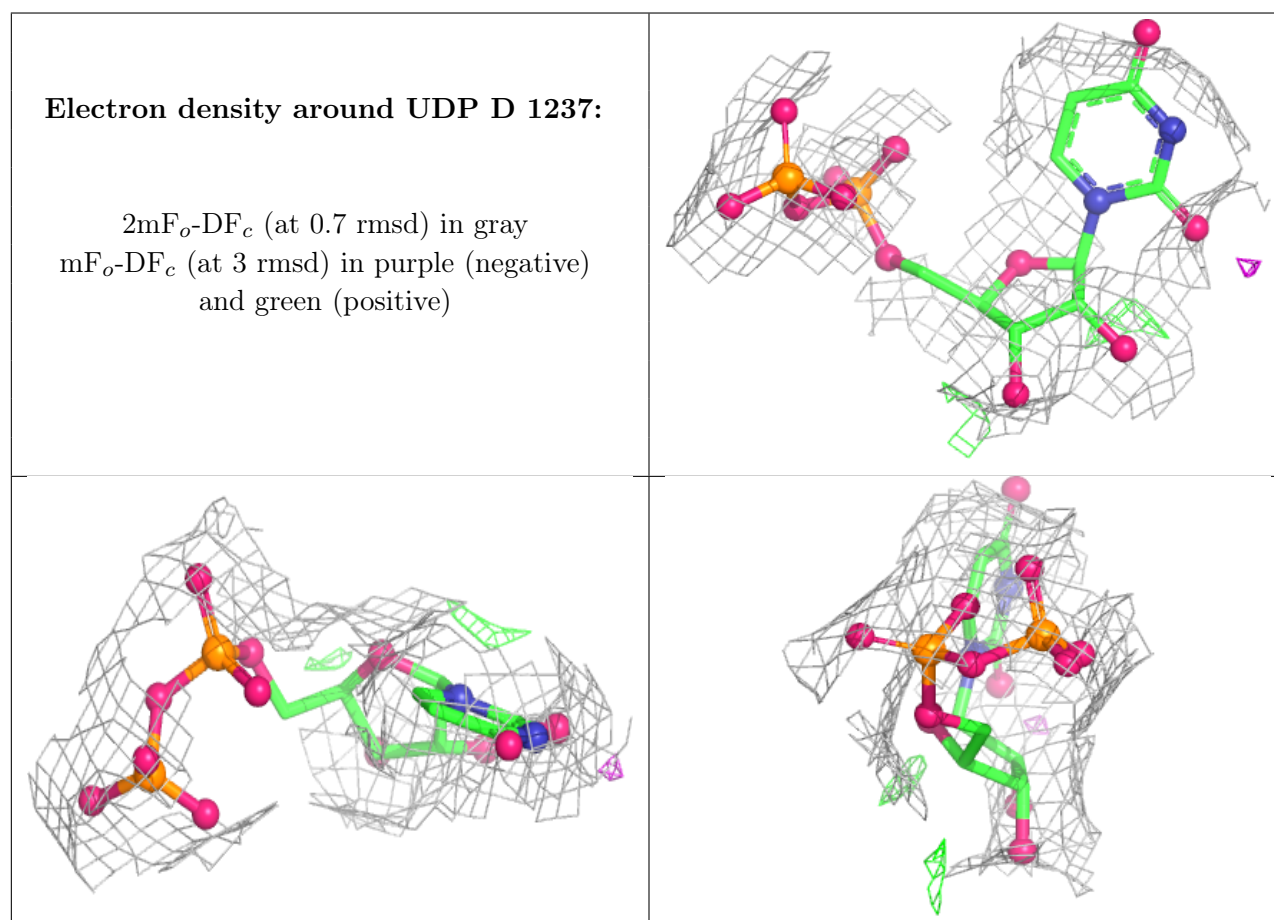
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

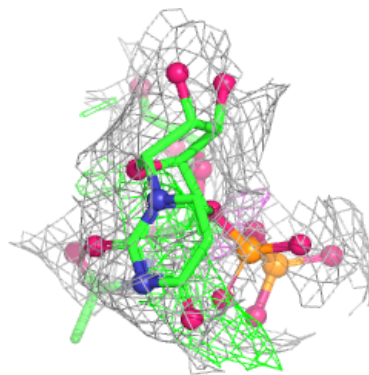
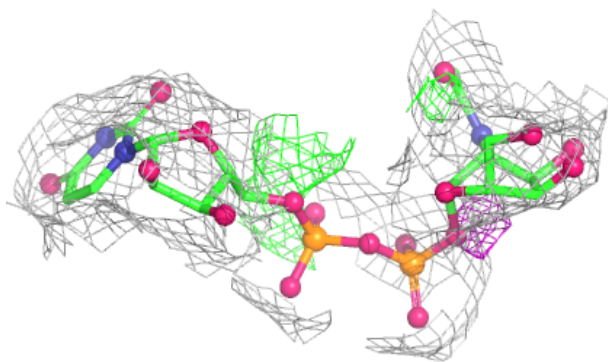
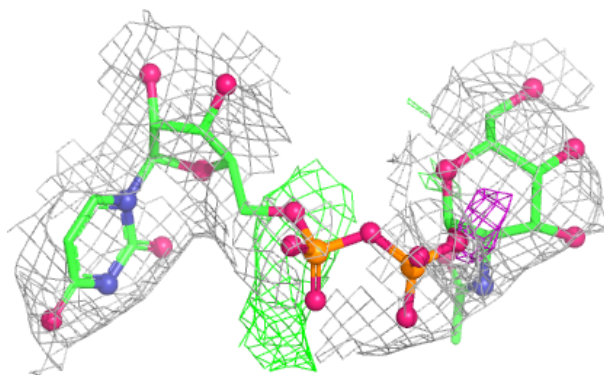
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	UDP	D	1237	25/25	0.67	0.12	97,127,136,141	0
4	A2G	D	1238	15/15	0.74	0.12	50,59,71,72	0
2	UD2	C	1241	39/39	0.79	0.12	69,86,97,98	0
3	UDP	B	1239	25/25	0.80	0.12	65,78,105,116	0
4	A2G	B	1240	15/15	0.82	0.14	62,79,86,89	0
2	UD2	A	1242	39/39	0.88	0.11	33,49,62,65	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.



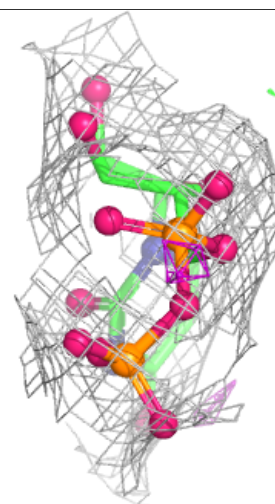
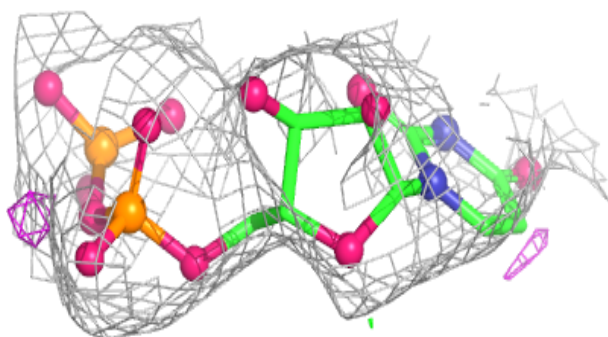
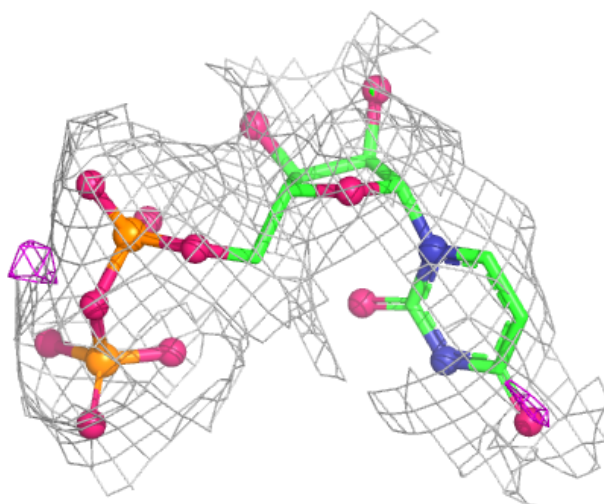
Electron density around UD2 C 1241:

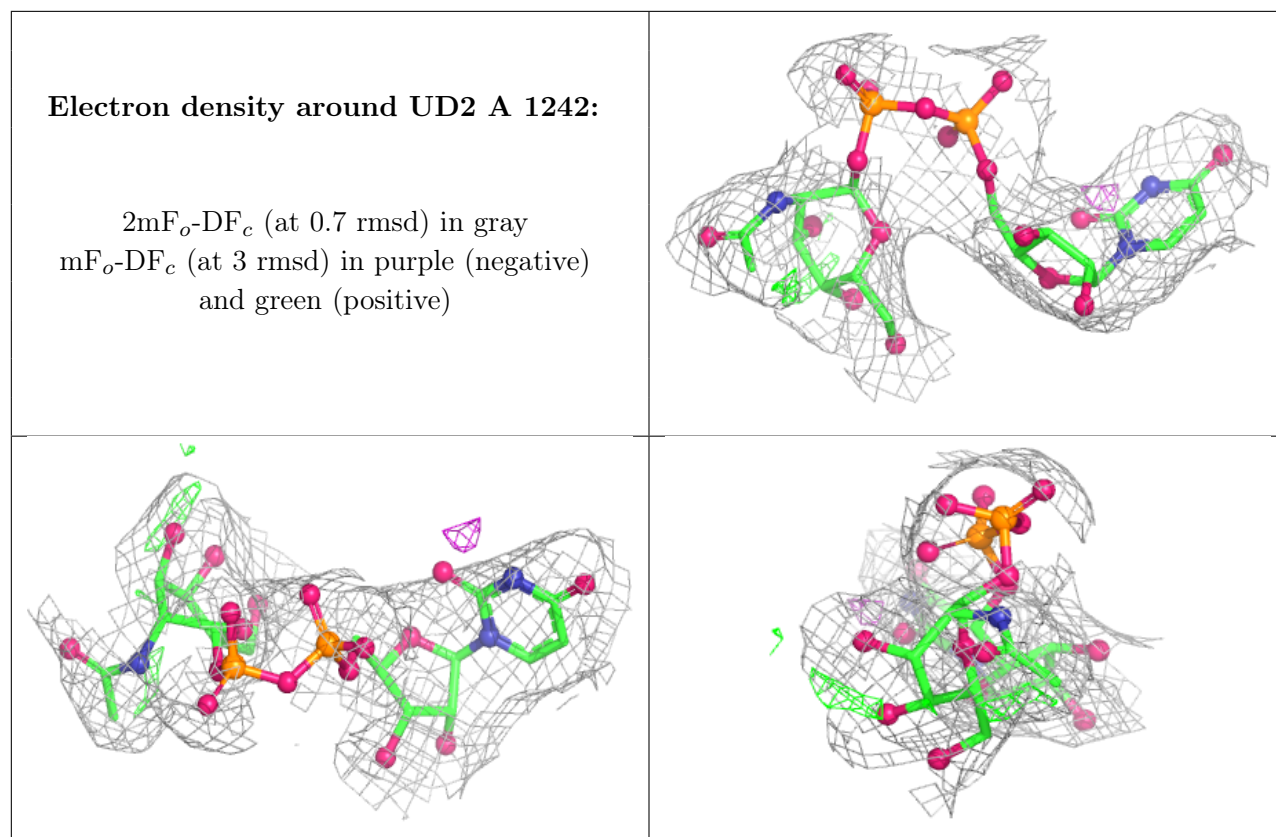
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around UDP B 1239:

$2mF_o-DF_c$ (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.