



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:14 AM UTC

PDB ID : 1B55 / pdb_00001b55
Title : PH DOMAIN FROM BRUTON'S TYROSINE KINASE IN COMPLEX WITH INOSITOL 1,3,4,5-TETRAKISPHOSPHATE
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Deposited on : 1999-01-12
Resolution : 2.40 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

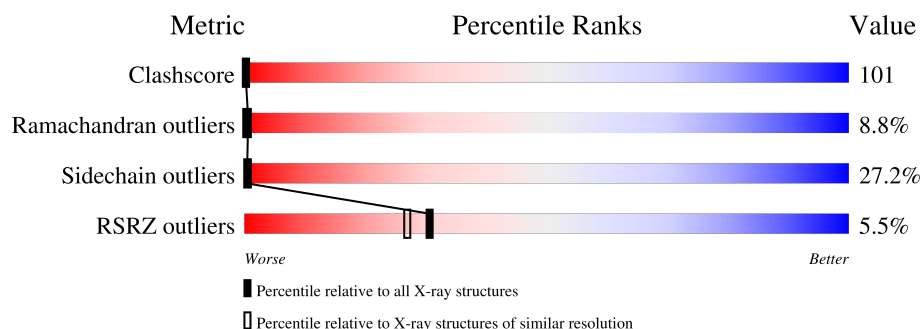
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	169	23% 46% 21% 6% .
1	B	169	9% 8% 37% 36% 15% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2964 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.

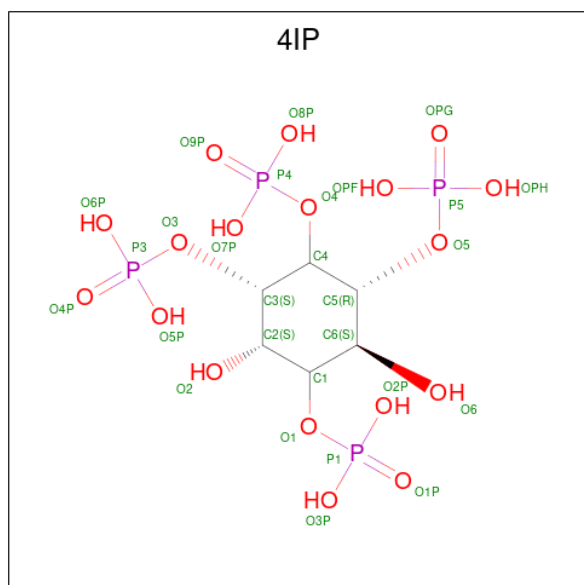
- Molecule 1 is a protein called TYROSINE-PROTEIN KINASE BTK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	163	Total	C	N	O	S	23	0	0
			1362	877	238	240	7			
1	B	162	Total	C	N	O	S	63	0	0
			1348	868	231	242	7			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is INOSITOL-(1,3,4,5)-TETRAKISPHOSPHATE (CCD ID: 4IP) (formula: C₆H₁₆O₁₈P₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			28	6	18	4		
3	B	1	Total	C	O	P	0	0
			28	6	18	4		

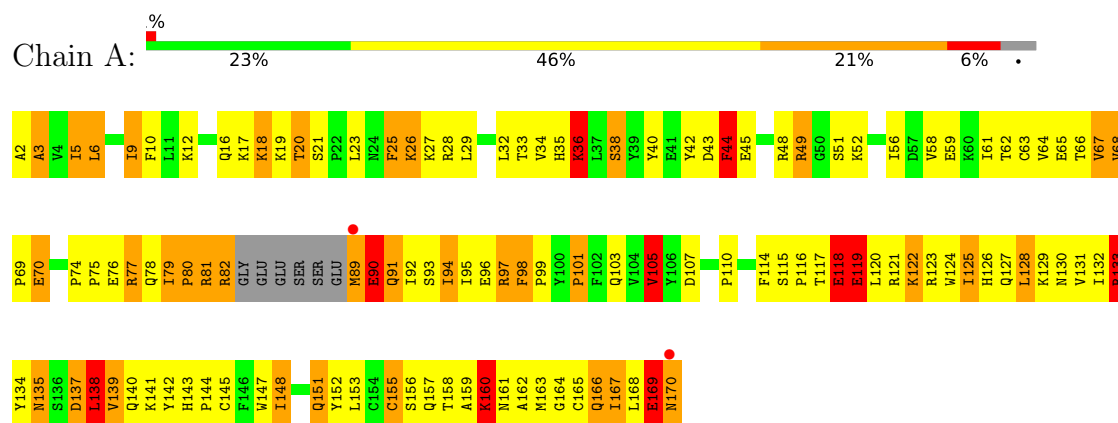
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	119	Total	O	0	0
			119	119		
4	B	77	Total	O	0	0
			77	77		

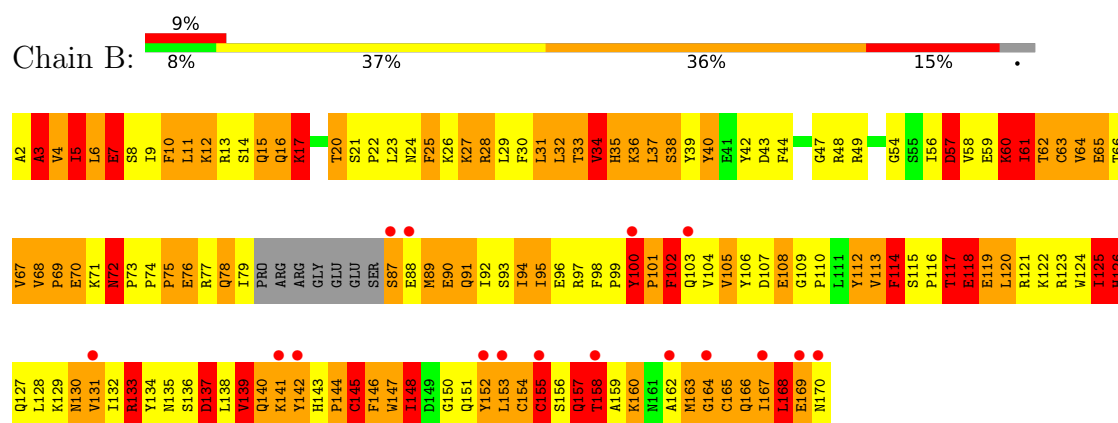
3 Residue-property plots

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: TYROSINE-PROTEIN KINASE BTK



• Molecule 1: TYROSINE-PROTEIN KINASE BTK



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	110.20Å 110.20Å 215.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.00 – 2.40 35.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.5 (35.00-2.40) 96.4 (35.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.39Å)	Xtriage
Refinement program	TNT 5E	Depositor
R, R_{free}	0.234 , 0.350 0.248 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 170.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2964	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: 4IP, ZN

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	1.58	16/1395 (1.1%)	1.75	21/1880 (1.1%)
1	B	1.58	13/1380 (0.9%)	2.05	68/1860 (3.7%)
All	All	1.58	29/2775 (1.0%)	1.91	89/3740 (2.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	1	5
All	All	1	9

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	125	ILE	CA-CB	-8.97	1.42	1.54
1	B	104	VAL	CA-CB	-7.11	1.45	1.54
1	A	138	LEU	N-CA	-7.05	1.37	1.46
1	A	56	ILE	CA-CB	-7.00	1.45	1.54
1	B	28	ARG	CA-C	-6.81	1.44	1.52
1	B	95	ILE	CA-C	-6.80	1.44	1.52
1	B	57	ASP	CA-C	-6.79	1.44	1.52
1	A	90	GLU	N-CA	-6.54	1.37	1.46
1	B	33	THR	CA-C	-6.47	1.46	1.53
1	A	99	PRO	N-CA	-6.23	1.42	1.47
1	B	117	THR	CA-C	5.97	1.59	1.52
1	A	90	GLU	CA-C	-5.88	1.45	1.53
1	B	100	TYR	CA-C	-5.70	1.45	1.52
1	B	10	PHE	CA-C	-5.63	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	7	GLU	CA-C	-5.55	1.45	1.52
1	A	6	LEU	CA-C	-5.46	1.46	1.52
1	A	101	PRO	CA-C	-5.43	1.47	1.52
1	A	38	SER	CA-C	-5.42	1.46	1.52
1	A	36	LYS	CA-C	-5.28	1.46	1.52
1	B	56	ILE	CA-C	-5.22	1.46	1.52
1	B	63	CYS	CA-C	5.18	1.58	1.52
1	B	4	VAL	C-O	-5.15	1.17	1.24
1	A	118	GLU	CA-C	-5.13	1.46	1.52
1	A	119	GLU	CA-C	-5.11	1.46	1.52
1	A	148	ILE	CA-CB	-5.11	1.49	1.55
1	A	167	ILE	CA-C	-5.11	1.46	1.52
1	A	26	LYS	C-O	-5.10	1.17	1.23
1	A	105	VAL	CA-CB	-5.06	1.47	1.54
1	A	95	ILE	CA-C	-5.00	1.46	1.52

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	148	ILE	N-CA-C	11.34	123.31	108.35
1	B	5	ILE	N-CA-C	-11.22	102.42	113.20
1	B	89	MET	N-CA-C	10.91	129.87	109.24
1	B	145	CYS	N-CA-C	10.76	123.64	110.41
1	A	168	LEU	CA-C-N	-8.09	111.37	122.77
1	A	168	LEU	C-N-CA	-8.09	111.37	122.77
1	B	104	VAL	CB-CA-C	-8.07	99.31	110.77
1	A	167	ILE	CB-CA-C	-7.81	99.55	110.96
1	B	72	ASN	O-C-N	7.52	127.32	121.23
1	B	100	TYR	O-C-N	7.34	129.76	121.32
1	B	61	ILE	N-CA-CB	7.21	118.39	110.53
1	B	155	CYS	N-CA-C	-7.09	97.34	108.90
1	B	71	LYS	N-CA-C	7.07	118.78	111.14
1	B	64	VAL	CA-C-N	-7.07	108.68	121.62
1	B	64	VAL	C-N-CA	-7.07	108.68	121.62
1	B	113	VAL	CB-CA-C	-7.00	102.22	110.91
1	B	101	PRO	CB-CA-C	-6.72	104.23	111.56
1	A	21	SER	N-CA-C	6.69	118.23	109.93
1	B	117	THR	N-CA-C	6.65	119.53	109.23
1	B	60	LYS	CA-C-N	-6.64	112.50	121.66
1	B	60	LYS	C-N-CA	-6.64	112.50	121.66
1	B	167	ILE	N-CA-C	-6.58	98.16	108.81
1	A	105	VAL	CB-CA-C	-6.55	100.40	110.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	118	GLU	CA-C-N	-6.42	109.28	121.54
1	B	118	GLU	C-N-CA	-6.42	109.28	121.54
1	B	10	PHE	N-CA-C	-6.25	100.95	110.14
1	B	22	PRO	CA-C-N	-6.20	114.16	122.72
1	B	22	PRO	C-N-CA	-6.20	114.16	122.72
1	A	61	ILE	N-CA-C	6.13	117.24	107.98
1	B	154	CYS	N-CA-C	-6.00	102.38	111.56
1	B	145	CYS	CB-CA-C	6.00	123.26	110.19
1	B	158	THR	N-CA-C	5.95	121.98	114.31
1	B	147	TRP	CA-C-N	-5.94	112.93	122.25
1	B	147	TRP	C-N-CA	-5.94	112.93	122.25
1	B	90	GLU	N-CA-C	5.91	118.39	107.99
1	B	27	LYS	CA-C-N	-5.88	112.68	122.21
1	B	27	LYS	C-N-CA	-5.88	112.68	122.21
1	A	44	PHE	N-CA-C	5.88	117.77	111.36
1	A	5	ILE	N-CA-C	-5.85	105.38	111.58
1	B	137	ASP	N-CA-C	5.73	117.17	110.41
1	B	27	LYS	N-CA-C	-5.73	99.95	109.46
1	B	101	PRO	CA-C-O	-5.72	114.22	120.92
1	A	121	ARG	N-CA-C	-5.72	105.12	111.36
1	B	38	SER	N-CA-C	5.72	118.87	109.72
1	B	102	PHE	CA-C-N	-5.70	111.97	121.89
1	B	102	PHE	C-N-CA	-5.70	111.97	121.89
1	B	5	ILE	CA-C-O	5.69	123.90	119.46
1	B	35	HIS	CA-CB-CG	-5.66	108.14	113.80
1	B	34	VAL	CA-C-N	-5.65	110.75	121.54
1	B	34	VAL	C-N-CA	-5.65	110.75	121.54
1	B	11	LEU	N-CA-C	-5.64	102.18	110.52
1	B	145	CYS	CA-CB-SG	-5.63	101.44	114.40
1	B	60	LYS	N-CA-C	-5.63	106.27	113.02
1	B	95	ILE	CB-CA-C	-5.57	102.15	111.29
1	B	20	THR	CB-CA-C	-5.55	102.16	110.88
1	A	67	VAL	N-CA-C	-5.54	99.82	108.86
1	B	95	ILE	N-CA-CB	5.54	120.37	111.23
1	B	164	GLY	N-CA-C	5.49	126.18	113.18
1	B	113	VAL	N-CA-C	5.48	116.47	108.58
1	B	87	SER	CA-C-N	5.48	129.02	120.75
1	B	87	SER	C-N-CA	5.48	129.02	120.75
1	B	151	GLN	N-CA-C	5.47	117.41	108.55
1	B	114	PHE	N-CA-C	5.46	118.63	109.95
1	B	112	TYR	N-CA-C	5.46	118.05	108.56
1	A	20	THR	N-CA-CB	-5.36	101.73	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	CYS	CA-C-O	5.31	125.13	119.34
1	A	165	CYS	N-CA-C	-5.29	106.41	112.92
1	B	141	LYS	N-CA-C	-5.27	100.44	109.24
1	B	90	GLU	CA-CB-CG	-5.25	103.59	114.10
1	B	3	ALA	N-CA-C	-5.22	99.68	110.80
1	A	98	PHE	CA-C-N	-5.21	114.74	121.04
1	A	98	PHE	C-N-CA	-5.21	114.74	121.04
1	B	34	VAL	N-CA-C	-5.21	106.72	111.67
1	B	168	LEU	N-CA-C	-5.20	99.94	108.41
1	B	31	LEU	N-CA-C	5.19	117.78	110.14
1	A	3	ALA	CA-C-N	-5.19	116.39	122.93
1	A	3	ALA	C-N-CA	-5.19	116.39	122.93
1	B	64	VAL	O-C-N	-5.17	116.56	122.62
1	B	35	HIS	CA-C-N	-5.17	113.60	123.32
1	B	35	HIS	C-N-CA	-5.17	113.60	123.32
1	A	18	LYS	N-CA-C	5.17	118.64	109.96
1	A	128	LEU	N-CA-C	-5.13	105.58	111.07
1	B	146	PHE	CA-C-N	-5.08	115.65	123.07
1	B	146	PHE	C-N-CA	-5.08	115.65	123.07
1	A	137	ASP	N-CA-C	5.08	119.03	111.87
1	B	125	ILE	N-CA-CB	-5.08	102.85	111.23
1	A	166	GLN	N-CA-C	5.07	117.67	109.06
1	B	67	VAL	N-CA-C	-5.04	98.85	109.34
1	B	56	ILE	N-CA-C	-5.02	101.39	108.27

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	145	CYS	CA

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	125	ILE	Mainchain
1	A	135	ASN	Sidechain
1	A	169	GLU	Mainchain
1	A	44	PHE	Sidechain
1	B	102	PHE	Sidechain
1	B	142	TYR	Sidechain
1	B	57	ASP	Sidechain
1	B	72	ASN	Sidechain
1	B	89	MET	Mainchain

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	1362	0	1378	136	3
1	B	1348	0	1358	411	0
2	A	1	0	0	0	0
2	B	1	0	0	1	0
3	A	28	0	8	2	0
3	B	28	0	8	2	0
4	A	119	0	0	13	4
4	B	77	0	0	58	1
All	All	2964	0	2752	538	7

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 101.

All (538) 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:B:101:PRO:HA	4:B:178:HOH:O	1.30	1.25
1:B:155:CYS:SG	1:B:165:CYS:HB3	1.79	1.22
1:A:97:ARG:HG3	1:A:97:ARG:HH11	1.06	1.17
1:B:100:TYR:HB3	1:B:121:ARG:HD3	1.31	1.11
1:B:87:SER:HB2	1:B:119:GLU:HB3	1.13	1.09
1:A:29:LEU:HD23	1:A:49:ARG:HB2	1.32	1.09
1:B:125:ILE:HD11	4:B:174:HOH:O	1.52	1.09
1:B:115:SER:CA	4:B:179:HOH:O	2.01	1.08
1:B:115:SER:HA	4:B:179:HOH:O	1.49	1.07
1:B:100:TYR:HB2	1:B:115:SER:HB3	1.38	1.05
1:B:100:TYR:CE1	1:B:118:GLU:HG3	1.93	1.02
1:B:105:VAL:HB	4:B:173:HOH:O	1.58	1.02
1:B:127:GLN:HA	4:B:181:HOH:O	1.58	1.02
1:B:135:ASN:HB2	1:B:138:LEU:HD21	1.03	1.02
1:B:40:TYR:CD2	1:B:49:ARG:HD2	1.96	0.99
1:B:10:PHE:CD1	4:B:179:HOH:O	2.13	0.99
1:B:115:SER:HB3	4:B:213:HOH:O	1.65	0.97
1:B:133:ARG:HG2	1:B:134:TYR:N	1.80	0.97
1:B:67:VAL:HG22	1:B:103:GLN:HB3	1.48	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:TYR:HB3	1:B:44:PHE:HE2	1.30	0.94
1:B:5:ILE:O	1:B:6:LEU:HD13	1.66	0.94
1:B:118:GLU:O	1:B:119:GLU:C	2.07	0.94
1:B:141:LYS:HA	1:B:167:ILE:HA	1.47	0.94
1:B:68:VAL:HB	4:B:230:HOH:O	1.66	0.94
1:B:143:HIS:HB3	1:B:154:CYS:CB	1.98	0.93
1:B:135:ASN:HB2	1:B:138:LEU:CD2	1.95	0.93
1:B:64:VAL:HG23	1:B:132:ILE:HD13	1.49	0.93
1:B:69:PRO:HG3	4:B:202:HOH:O	1.67	0.93
1:B:99:PRO:O	1:B:101:PRO:HD3	1.67	0.93
1:B:135:ASN:CB	1:B:138:LEU:HD21	1.97	0.92
1:B:90:GLU:HB3	1:B:93:SER:OG	1.69	0.92
1:B:67:VAL:CG2	1:B:103:GLN:HB3	1.98	0.92
1:B:125:ILE:HG22	1:B:126:HIS:N	1.82	0.92
1:B:13:ARG:HH12	1:B:148:ILE:HG23	1.33	0.92
1:A:62:THR:HB	1:A:139:VAL:HG13	1.50	0.91
1:B:30:PHE:C	1:B:31:LEU:HD23	1.96	0.90
1:A:59:GLU:HA	1:A:135:ASN:HD21	1.36	0.90
1:B:121:ARG:HD2	4:B:194:HOH:O	1.71	0.89
1:B:57:ASP:HB2	1:B:60:LYS:HG3	1.52	0.89
1:B:100:TYR:HE1	1:B:118:GLU:HG3	1.37	0.89
1:B:87:SER:CB	1:B:119:GLU:HB3	2.02	0.89
1:B:90:GLU:OE1	1:B:90:GLU:HA	1.70	0.88
1:A:2:ALA:HB1	4:A:258:HOH:O	1.73	0.88
1:A:118:GLU:HG3	4:A:216:HOH:O	1.74	0.88
1:B:64:VAL:HG11	1:B:128:LEU:CB	2.03	0.87
1:B:98:PHE:HA	1:B:116:PRO:HA	1.56	0.87
1:B:2:ALA:HB3	1:B:33:THR:CG2	2.05	0.87
1:A:58:VAL:HG11	1:A:131:VAL:HG12	1.56	0.86
1:A:89:MET:HA	4:A:173:HOH:O	1.73	0.86
1:B:13:ARG:NH2	1:B:112:TYR:HE2	1.73	0.86
1:A:42:TYR:HB3	1:B:44:PHE:CE2	2.10	0.86
1:B:13:ARG:NH2	1:B:112:TYR:CE2	2.44	0.85
1:B:73:PRO:O	1:B:78:GLN:HG2	1.76	0.85
1:B:61:ILE:HB	1:B:132:ILE:HG21	1.57	0.85
1:B:63:CYS:HB3	4:B:173:HOH:O	1.74	0.85
1:B:147:TRP:C	1:B:148:ILE:HG13	2.02	0.84
1:B:102:PHE:HB3	4:B:194:HOH:O	1.77	0.84
1:B:143:HIS:CE1	1:B:155:CYS:SG	2.71	0.84
1:B:67:VAL:HG23	1:B:102:PHE:HA	1.59	0.84
1:B:130:ASN:HB2	4:B:181:HOH:O	1.77	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:HIS:CE1	1:B:164:GLY:HA2	2.13	0.83
1:B:32:LEU:HD12	1:B:33:THR:N	1.92	0.82
1:B:88:GLU:HB2	1:B:117:THR:CG2	2.09	0.82
1:B:143:HIS:HB3	1:B:154:CYS:HB3	1.60	0.82
1:A:92:ILE:HD11	1:B:9:ILE:HG13	1.62	0.82
1:A:34:VAL:HG23	1:A:35:HIS:HD2	1.43	0.82
1:B:5:ILE:O	1:B:6:LEU:CD1	2.27	0.82
1:B:92:ILE:O	1:B:96:GLU:HG3	1.80	0.82
1:B:13:ARG:HH21	1:B:112:TYR:HE2	1.25	0.82
1:B:168:LEU:N	4:B:177:HOH:O	2.13	0.81
1:A:97:ARG:HG3	1:A:97:ARG:NH1	1.86	0.81
1:B:155:CYS:HB2	1:B:157:GLN:HG3	1.60	0.81
1:B:90:GLU:O	1:B:92:ILE:N	2.13	0.81
1:A:45:GLU:O	1:B:27:LYS:HE3	1.81	0.81
1:A:90:GLU:O	1:A:92:ILE:N	2.14	0.81
1:B:159:ALA:O	1:B:160:LYS:C	2.22	0.81
1:B:75:PRO:CA	1:B:78:GLN:HG3	2.11	0.81
1:B:64:VAL:HG21	1:B:128:LEU:HB3	1.63	0.80
1:A:5:ILE:HD12	1:A:32:LEU:O	1.82	0.79
1:B:114:PHE:C	4:B:179:HOH:O	2.24	0.79
1:B:77:ARG:O	1:B:99:PRO:HB3	1.81	0.79
1:B:68:VAL:O	1:B:69:PRO:C	2.25	0.79
1:B:127:GLN:O	1:B:131:VAL:HG22	1.83	0.79
1:A:81:ARG:HA	4:A:271:HOH:O	1.83	0.79
1:B:40:TYR:CE2	1:B:49:ARG:HD2	2.17	0.78
1:B:90:GLU:HB3	1:B:93:SER:CB	2.13	0.78
1:B:100:TYR:HB3	1:B:121:ARG:CD	2.11	0.78
1:A:58:VAL:HG11	1:A:131:VAL:CG1	2.14	0.78
1:B:6:LEU:HB3	1:B:124:TRP:CZ2	2.19	0.78
1:B:74:PRO:O	1:B:75:PRO:C	2.27	0.78
1:B:99:PRO:HD2	1:B:100:TYR:CD2	2.18	0.77
1:B:67:VAL:HG11	1:B:146:PHE:CZ	2.19	0.77
1:B:63:CYS:N	4:B:173:HOH:O	2.19	0.76
1:A:16:GLN:OE1	1:A:23:LEU:HD13	1.84	0.76
1:B:126:HIS:HD2	4:B:204:HOH:O	1.67	0.76
1:B:100:TYR:CD1	1:B:118:GLU:HG3	2.19	0.76
1:B:115:SER:CB	4:B:213:HOH:O	2.28	0.76
1:A:80:PRO:HG3	1:A:97:ARG:HE	1.49	0.75
1:A:91:GLN:O	1:A:91:GLN:HG2	1.84	0.75
1:B:31:LEU:HD23	1:B:31:LEU:N	1.98	0.75
1:B:25:PHE:CD1	1:B:25:PHE:N	2.53	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:VAL:HG12	1:B:6:LEU:H	1.50	0.75
1:B:140:GLN:O	1:B:167:ILE:HG23	1.86	0.74
1:B:75:PRO:O	1:B:78:GLN:N	2.20	0.73
1:B:145:CYS:HA	4:B:230:HOH:O	1.86	0.73
1:B:146:PHE:C	1:B:148:ILE:HD11	2.13	0.73
1:B:143:HIS:CE1	1:B:165:CYS:H	2.05	0.73
1:B:166:GLN:HG3	1:B:167:ILE:H	1.52	0.73
1:B:64:VAL:HG23	1:B:132:ILE:CD1	2.19	0.73
1:B:125:ILE:O	1:B:128:LEU:HB2	1.89	0.73
1:B:35:HIS:HB3	4:B:228:HOH:O	1.88	0.72
1:B:8:SER:O	1:B:29:LEU:HD12	1.89	0.72
1:B:121:ARG:HB3	4:B:192:HOH:O	1.89	0.72
1:A:81:ARG:O	1:A:82:ARG:HD2	1.90	0.72
1:A:101:PRO:HB3	1:A:114:PHE:CD2	2.24	0.72
1:B:101:PRO:HD2	4:B:202:HOH:O	1.89	0.72
1:B:62:THR:HG23	1:B:106:TYR:HA	1.72	0.72
1:B:107:ASP:OD2	1:B:108:GLU:N	2.23	0.71
1:B:64:VAL:HG11	1:B:128:LEU:HB2	1.70	0.71
1:A:58:VAL:CG1	1:A:131:VAL:HG12	2.21	0.71
1:A:97:ARG:HH11	1:A:97:ARG:CG	1.96	0.70
1:B:30:PHE:O	1:B:31:LEU:HD23	1.92	0.70
1:B:67:VAL:CG2	1:B:102:PHE:C	2.64	0.70
1:B:125:ILE:O	1:B:126:HIS:C	2.31	0.70
1:B:64:VAL:HG11	1:B:128:LEU:HB3	1.73	0.70
1:B:67:VAL:HG12	1:B:68:VAL:H	1.57	0.70
1:B:114:PHE:HD1	1:B:114:PHE:N	1.90	0.70
1:B:114:PHE:N	1:B:114:PHE:CD1	2.60	0.69
1:B:31:LEU:O	1:B:37:LEU:HD12	1.93	0.69
1:B:64:VAL:HG13	1:B:128:LEU:HD13	1.74	0.69
1:B:143:HIS:O	1:B:145:CYS:N	2.26	0.69
1:B:13:ARG:NH1	1:B:148:ILE:HG23	2.07	0.69
1:B:5:ILE:HD12	1:B:33:THR:HA	1.73	0.69
1:B:87:SER:HB2	1:B:119:GLU:CB	2.07	0.69
1:B:100:TYR:HB2	1:B:115:SER:CB	2.20	0.69
3:A:171:4IP:O5P	4:A:197:HOH:O	2.10	0.69
1:B:165:CYS:O	1:B:166:GLN:HB2	1.93	0.68
1:B:79:ILE:HG22	1:B:79:ILE:O	1.94	0.68
1:B:141:LYS:HB3	1:B:166:GLN:O	1.94	0.68
1:A:141:LYS:NZ	1:A:163:MET:HE1	2.09	0.67
1:B:118:GLU:O	1:B:120:LEU:N	2.27	0.67
1:A:34:VAL:HG23	1:A:35:HIS:CD2	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:CYS:SG	1:A:157:GLN:HG3	2.34	0.67
1:B:88:GLU:HB2	1:B:117:THR:HB	1.76	0.67
1:A:81:ARG:HG2	4:A:271:HOH:O	1.93	0.67
1:B:58:VAL:O	1:B:61:ILE:HG13	1.94	0.67
1:B:32:LEU:HD12	1:B:33:THR:H	1.59	0.67
1:B:147:TRP:C	1:B:148:ILE:CG1	2.68	0.67
1:A:137:ASP:O	1:A:139:VAL:N	2.28	0.66
1:B:61:ILE:HB	1:B:132:ILE:CG2	2.25	0.66
1:B:90:GLU:O	1:B:93:SER:N	2.28	0.66
1:B:166:GLN:HG3	1:B:167:ILE:N	2.09	0.66
1:B:13:ARG:HD2	1:B:23:LEU:HD21	1.75	0.66
1:B:88:GLU:HB2	1:B:117:THR:CB	2.26	0.66
1:A:143:HIS:CE1	1:A:164:GLY:HA2	2.30	0.66
1:B:145:CYS:CA	4:B:230:HOH:O	2.42	0.65
1:B:129:LYS:O	1:B:130:ASN:C	2.38	0.65
1:B:12:LYS:O	1:B:25:PHE:HA	1.96	0.65
1:B:129:LYS:O	1:B:131:VAL:N	2.29	0.65
1:B:142:TYR:OH	1:B:144:PRO:HB3	1.96	0.65
1:B:7:GLU:HA	1:B:30:PHE:O	1.95	0.65
1:A:147:TRP:HD1	1:A:152:TYR:CE2	2.14	0.65
1:B:4:VAL:CG1	1:B:6:LEU:O	2.45	0.65
1:B:75:PRO:C	1:B:78:GLN:HG3	2.22	0.65
1:B:127:GLN:CA	4:B:181:HOH:O	2.30	0.65
1:B:2:ALA:HB3	1:B:33:THR:HG21	1.78	0.65
1:B:155:CYS:HB2	1:B:157:GLN:CG	2.27	0.65
1:A:63:CYS:O	1:A:105:VAL:HG23	1.96	0.65
1:B:100:TYR:CD2	1:B:100:TYR:N	2.64	0.64
1:B:159:ALA:O	1:B:162:ALA:N	2.26	0.64
1:A:10:PHE:CE1	1:A:115:SER:HB2	2.33	0.64
1:B:90:GLU:OE1	1:B:90:GLU:CA	2.37	0.64
1:B:67:VAL:HG23	1:B:102:PHE:CA	2.27	0.64
1:B:67:VAL:HG11	1:B:146:PHE:CE1	2.33	0.64
1:A:80:PRO:CG	1:A:97:ARG:HE	2.11	0.64
1:A:29:LEU:CD2	1:A:49:ARG:HB2	2.19	0.63
1:B:99:PRO:HD2	1:B:100:TYR:CE2	2.32	0.63
1:B:99:PRO:HD2	1:B:100:TYR:HD2	1.61	0.63
1:B:2:ALA:HB3	1:B:33:THR:HG22	1.81	0.63
1:B:68:VAL:O	1:B:68:VAL:HG12	1.99	0.63
1:A:119:GLU:HG2	1:A:119:GLU:O	1.91	0.63
1:B:20:THR:OG1	1:B:21:SER:N	2.28	0.63
1:A:70:GLU:O	1:A:70:GLU:HG2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:HG21	1:B:103:GLN:HB3	1.81	0.62
1:A:5:ILE:HD12	1:A:33:THR:HA	1.81	0.62
1:A:159:ALA:O	1:A:162:ALA:N	2.30	0.62
1:B:94:ILE:HA	4:B:188:HOH:O	1.99	0.62
1:B:154:CYS:SG	1:B:155:CYS:N	2.73	0.62
1:B:72:ASN:ND2	4:B:224:HOH:O	2.33	0.62
1:B:99:PRO:HG2	1:B:100:TYR:CE2	2.34	0.62
1:A:141:LYS:HE3	1:A:163:MET:HE2	1.82	0.61
1:B:169:GLU:HG2	1:B:170:ASN:N	2.11	0.61
1:A:101:PRO:HB3	1:A:114:PHE:CE2	2.34	0.61
1:A:43:ASP:HB3	1:A:48:ARG:O	2.00	0.61
1:A:29:LEU:HD23	1:A:49:ARG:CB	2.21	0.61
1:A:141:LYS:HZ2	1:A:163:MET:HE1	1.65	0.61
1:B:67:VAL:HG21	1:B:102:PHE:O	2.01	0.61
1:B:165:CYS:O	1:B:165:CYS:SG	2.59	0.61
1:A:6:LEU:HB3	1:A:124:TRP:CZ2	2.36	0.61
1:B:98:PHE:HA	1:B:115:SER:O	2.01	0.61
1:B:154:CYS:C	1:B:165:CYS:SG	2.84	0.61
1:B:121:ARG:NH1	4:B:202:HOH:O	2.23	0.61
1:B:147:TRP:CZ2	1:B:150:GLY:O	2.54	0.61
1:A:59:GLU:CA	1:A:135:ASN:HD21	2.13	0.60
1:A:90:GLU:O	1:A:91:GLN:C	2.43	0.60
1:B:58:VAL:HA	1:B:61:ILE:HG13	1.83	0.60
1:B:121:ARG:O	1:B:122:LYS:C	2.45	0.60
1:A:89:MET:HB3	4:A:219:HOH:O	2.00	0.60
1:A:94:ILE:HD13	1:B:91:GLN:HG3	1.83	0.60
1:B:98:PHE:CA	1:B:115:SER:O	2.49	0.60
1:A:166:GLN:HG2	1:A:167:ILE:N	2.15	0.60
1:B:115:SER:N	4:B:179:HOH:O	2.25	0.60
1:B:92:ILE:HB	4:B:219:HOH:O	2.01	0.60
1:B:4:VAL:HG12	1:B:6:LEU:N	2.16	0.60
1:B:154:CYS:O	1:B:165:CYS:SG	2.59	0.60
1:B:155:CYS:SG	1:B:165:CYS:CB	2.72	0.60
1:B:113:VAL:C	4:B:178:HOH:O	2.43	0.60
1:B:128:LEU:O	1:B:131:VAL:HG23	2.02	0.59
1:B:75:PRO:O	1:B:78:GLN:HG3	2.02	0.59
1:B:75:PRO:O	1:B:77:ARG:N	2.35	0.59
1:B:21:SER:OG	3:B:172:4IP:OPH	2.20	0.59
1:B:67:VAL:CG1	1:B:146:PHE:CZ	2.86	0.59
1:B:40:TYR:CD1	1:B:40:TYR:N	2.71	0.59
1:A:5:ILE:HD12	1:A:33:THR:CA	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:HG23	1:B:102:PHE:C	2.27	0.59
1:B:4:VAL:CG1	1:B:6:LEU:H	2.14	0.59
1:A:141:LYS:HZ2	1:A:163:MET:CE	2.15	0.59
1:B:14:SER:N	1:B:24:ASN:O	2.35	0.59
1:B:148:ILE:N	1:B:148:ILE:HD12	2.17	0.59
1:B:67:VAL:HG12	1:B:146:PHE:CE2	2.38	0.58
1:B:73:PRO:O	1:B:78:GLN:NE2	2.35	0.58
1:A:145:CYS:HB3	1:A:153:LEU:HD12	1.86	0.58
1:B:143:HIS:ND1	1:B:154:CYS:SG	2.76	0.58
1:A:129:LYS:O	1:A:130:ASN:C	2.44	0.58
1:B:42:TYR:CE2	1:B:44:PHE:HA	2.38	0.58
1:A:76:GLU:HG3	1:A:77:ARG:HG2	1.86	0.58
1:B:26:LYS:HB3	1:B:28:ARG:CZ	2.32	0.58
1:B:98:PHE:HA	1:B:116:PRO:CA	2.31	0.58
1:A:25:PHE:CD1	1:A:25:PHE:N	2.70	0.58
1:B:2:ALA:O	1:B:3:ALA:HB2	2.04	0.58
1:B:34:VAL:O	1:B:58:VAL:HG21	2.03	0.58
1:B:40:TYR:CD2	1:B:49:ARG:CD	2.81	0.58
1:B:97:ARG:O	1:B:116:PRO:HA	2.03	0.58
1:B:67:VAL:HG21	1:B:102:PHE:C	2.28	0.58
1:B:125:ILE:O	1:B:128:LEU:N	2.37	0.58
1:A:64:VAL:C	1:A:65:GLU:HG2	2.28	0.57
1:B:13:ARG:HH11	1:B:23:LEU:HD21	1.69	0.57
1:B:143:HIS:C	1:B:145:CYS:N	2.61	0.57
1:B:62:THR:OG1	4:B:173:HOH:O	2.17	0.57
1:B:112:TYR:CZ	1:B:146:PHE:CD1	2.93	0.57
1:B:43:ASP:N	1:B:48:ARG:O	2.27	0.57
1:B:36:LYS:HA	1:B:58:VAL:HG23	1.86	0.57
1:B:49:ARG:HG2	1:B:49:ARG:HH11	1.68	0.57
1:B:115:SER:CB	4:B:179:HOH:O	2.41	0.57
1:B:92:ILE:N	4:B:219:HOH:O	2.27	0.57
1:B:90:GLU:O	1:B:91:GLN:C	2.47	0.56
1:A:98:PHE:HA	1:A:116:PRO:HA	1.86	0.56
1:A:115:SER:OG	1:A:116:PRO:HD2	2.05	0.56
1:B:75:PRO:N	1:B:78:GLN:HG3	2.19	0.56
1:B:74:PRO:C	1:B:78:GLN:HG3	2.30	0.56
1:B:35:HIS:O	1:B:58:VAL:HG23	2.04	0.56
1:B:139:VAL:C	1:B:140:GLN:HG2	2.31	0.56
1:A:92:ILE:HD11	1:B:9:ILE:CG1	2.34	0.56
1:B:64:VAL:O	1:B:65:GLU:HG2	2.06	0.56
1:B:143:HIS:CG	1:B:154:CYS:SG	2.98	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:GLN:HB3	1:A:158:THR:HG22	1.88	0.56
1:B:67:VAL:CG1	1:B:146:PHE:CE2	2.89	0.56
1:B:117:THR:N	4:B:213:HOH:O	2.39	0.56
1:B:126:HIS:HE1	4:B:209:HOH:O	1.88	0.56
1:B:99:PRO:HG2	1:B:100:TYR:HE2	1.70	0.56
1:A:5:ILE:HD12	1:A:32:LEU:C	2.30	0.55
1:A:58:VAL:HG12	1:A:58:VAL:O	2.07	0.55
1:B:150:GLY:O	1:B:158:THR:HG23	2.05	0.55
1:B:114:PHE:HA	4:B:178:HOH:O	2.07	0.55
1:B:4:VAL:HG12	1:B:5:ILE:N	2.21	0.55
1:B:7:GLU:N	1:B:124:TRP:HZ2	2.05	0.55
1:B:146:PHE:HB3	1:B:148:ILE:CG1	2.37	0.55
1:B:119:GLU:CD	1:B:123:ARG:HD2	2.32	0.55
1:B:147:TRP:CH2	1:B:150:GLY:C	2.85	0.55
1:A:3:ALA:O	1:A:33:THR:HG22	2.06	0.54
1:B:141:LYS:CA	1:B:166:GLN:O	2.55	0.54
1:A:134:TYR:HA	4:A:203:HOH:O	2.07	0.54
1:A:152:TYR:O	1:A:156:SER:N	2.39	0.54
1:B:13:ARG:NE	4:B:184:HOH:O	2.41	0.54
1:B:34:VAL:HG12	1:B:131:VAL:HG12	1.89	0.54
1:B:39:TYR:CZ	1:B:54:GLY:HA3	2.42	0.54
1:B:77:ARG:O	1:B:78:GLN:O	2.25	0.54
1:B:78:GLN:O	1:B:99:PRO:HB3	2.07	0.54
1:B:124:TRP:O	1:B:125:ILE:C	2.49	0.54
1:B:47:GLY:O	1:B:48:ARG:HB3	2.06	0.54
1:B:142:TYR:CE1	1:B:144:PRO:N	2.76	0.54
1:B:25:PHE:N	1:B:25:PHE:HD1	2.05	0.53
1:B:64:VAL:HG13	1:B:128:LEU:CD1	2.37	0.53
1:B:112:TYR:CZ	1:B:146:PHE:HD1	2.26	0.53
3:B:172:4IP:O9P	3:B:172:4IP:OPF	2.26	0.53
1:A:128:LEU:O	1:A:132:ILE:HG12	2.09	0.53
1:A:5:ILE:CD1	1:A:33:THR:HA	2.39	0.53
1:B:141:LYS:CB	1:B:166:GLN:O	2.56	0.53
1:A:169:GLU:C	1:A:170:ASN:HD22	2.17	0.53
1:B:143:HIS:CE1	1:B:164:GLY:CA	2.88	0.53
1:B:143:HIS:O	1:B:144:PRO:C	2.52	0.53
1:B:146:PHE:C	1:B:148:ILE:CD1	2.80	0.53
1:A:159:ALA:O	1:A:160:LYS:C	2.52	0.52
1:B:145:CYS:SG	1:B:154:CYS:HA	2.49	0.52
1:B:143:HIS:ND1	1:B:155:CYS:SG	2.82	0.52
1:B:64:VAL:CG2	1:B:132:ILE:HD13	2.31	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:TYR:CD2	1:A:52:LYS:HA	2.45	0.52
1:B:147:TRP:N	1:B:148:ILE:HG13	2.25	0.52
1:A:114:PHE:CD1	1:A:114:PHE:N	2.77	0.52
1:B:10:PHE:HD1	4:B:179:HOH:O	1.74	0.52
1:B:90:GLU:O	1:B:90:GLU:HG3	2.06	0.52
1:B:100:TYR:CB	1:B:121:ARG:HD3	2.22	0.52
1:B:159:ALA:O	1:B:162:ALA:CB	2.58	0.52
1:B:95:ILE:O	1:B:95:ILE:HG13	2.09	0.52
1:A:125:ILE:O	1:A:129:LYS:HG3	2.09	0.52
1:B:7:GLU:O	1:B:8:SER:OG	2.27	0.52
1:B:75:PRO:CA	1:B:78:GLN:CG	2.86	0.52
1:A:62:THR:O	1:A:138:LEU:HA	2.10	0.51
1:A:131:VAL:HG23	4:A:221:HOH:O	2.10	0.51
1:B:5:ILE:N	1:B:32:LEU:O	2.43	0.51
1:B:20:THR:O	1:B:21:SER:C	2.52	0.51
1:A:141:LYS:NZ	1:A:163:MET:CE	2.73	0.51
1:B:135:ASN:HB3	4:B:208:HOH:O	2.10	0.51
1:B:137:ASP:O	1:B:139:VAL:N	2.44	0.51
1:B:64:VAL:CG2	1:B:128:LEU:HB3	2.38	0.51
1:B:142:TYR:CE1	1:B:144:PRO:CA	2.93	0.51
1:B:143:HIS:HB3	1:B:154:CYS:HB2	1.88	0.51
1:B:127:GLN:HA	1:B:127:GLN:OE1	2.11	0.51
1:B:133:ARG:HG2	1:B:134:TYR:CD1	2.46	0.51
1:A:92:ILE:CD1	1:B:9:ILE:CD1	2.89	0.51
1:B:10:PHE:HZ	1:B:124:TRP:CD1	2.29	0.51
1:B:102:PHE:C	4:B:244:HOH:O	2.53	0.51
1:B:143:HIS:HE1	1:B:155:CYS:SG	2.29	0.51
1:B:75:PRO:HA	1:B:78:GLN:CD	2.36	0.50
1:A:137:ASP:O	1:A:138:LEU:C	2.52	0.50
1:B:90:GLU:HB3	1:B:93:SER:HB3	1.91	0.50
1:A:58:VAL:CG1	1:A:58:VAL:O	2.59	0.50
1:B:69:PRO:HD3	4:B:202:HOH:O	2.12	0.50
1:B:90:GLU:HG3	1:B:92:ILE:H	1.75	0.50
1:B:75:PRO:O	1:B:76:GLU:C	2.54	0.50
1:B:90:GLU:CG	1:B:93:SER:H	2.24	0.50
1:A:129:LYS:O	1:A:132:ILE:N	2.45	0.50
1:B:68:VAL:O	1:B:70:GLU:N	2.44	0.50
1:B:74:PRO:O	1:B:77:ARG:N	2.32	0.50
1:B:90:GLU:HG3	1:B:93:SER:H	1.75	0.50
1:A:63:CYS:HB2	1:A:139:VAL:HG22	1.93	0.50
3:A:171:4IP:O9P	3:A:171:4IP:OPF	2.30	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:PRO:CA	4:B:178:HOH:O	2.14	0.50
1:A:62:THR:HB	1:A:139:VAL:CG1	2.31	0.50
1:B:128:LEU:O	1:B:131:VAL:N	2.39	0.50
1:A:138:LEU:N	1:A:138:LEU:HD23	2.23	0.49
1:A:141:LYS:HE3	1:A:163:MET:CE	2.40	0.49
1:B:113:VAL:O	4:B:178:HOH:O	2.20	0.49
1:B:7:GLU:C	1:B:8:SER:OG	2.55	0.49
1:B:34:VAL:O	1:B:35:HIS:C	2.49	0.49
1:A:159:ALA:O	1:A:161:ASN:N	2.46	0.49
1:B:155:CYS:CB	1:B:165:CYS:HB3	2.42	0.49
1:B:118:GLU:C	1:B:120:LEU:N	2.68	0.49
1:B:74:PRO:O	1:B:78:GLN:HG3	2.13	0.49
1:B:132:ILE:O	1:B:134:TYR:N	2.46	0.49
1:A:169:GLU:O	1:A:170:ASN:C	2.56	0.49
1:B:13:ARG:HE	1:B:112:TYR:HD2	1.60	0.49
1:B:152:TYR:C	1:B:154:CYS:H	2.20	0.49
1:A:90:GLU:HA	1:A:93:SER:OG	2.12	0.49
1:B:4:VAL:CG1	1:B:5:ILE:N	2.76	0.49
1:A:34:VAL:CG2	1:A:35:HIS:HD2	2.21	0.49
1:B:7:GLU:HG3	1:B:31:LEU:CD2	2.43	0.49
1:B:30:PHE:CZ	1:B:113:VAL:HG22	2.48	0.49
1:A:42:TYR:HD2	1:B:44:PHE:CE2	2.31	0.48
1:B:57:ASP:HB2	1:B:60:LYS:CG	2.33	0.48
1:A:76:GLU:HG2	4:A:234:HOH:O	2.13	0.48
1:B:133:ARG:CG	1:B:134:TYR:CD1	2.97	0.48
1:A:80:PRO:HB3	1:A:97:ARG:HH21	1.77	0.48
1:B:77:ARG:HD2	1:B:114:PHE:CE2	2.48	0.48
1:B:90:GLU:HG3	4:B:219:HOH:O	2.13	0.48
1:B:99:PRO:N	1:B:115:SER:O	2.47	0.48
1:B:119:GLU:O	1:B:119:GLU:HG2	2.13	0.48
1:B:141:LYS:HA	1:B:166:GLN:O	2.13	0.48
1:B:35:HIS:O	1:B:58:VAL:CG2	2.62	0.48
1:B:63:CYS:CA	4:B:173:HOH:O	2.61	0.48
1:B:106:TYR:N	1:B:109:GLY:O	2.35	0.48
1:B:143:HIS:CE1	1:B:165:CYS:N	2.79	0.48
1:B:117:THR:C	4:B:213:HOH:O	2.55	0.48
1:B:15:GLN:HB3	1:B:17:LYS:HD2	1.95	0.48
1:B:94:ILE:O	1:B:94:ILE:HG23	2.13	0.48
1:B:10:PHE:HB3	1:B:114:PHE:O	2.14	0.47
1:A:10:PHE:HD1	1:A:115:SER:HA	1.79	0.47
1:B:88:GLU:HB2	1:B:117:THR:HG21	1.92	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:TYR:CZ	1:B:144:PRO:HG3	2.49	0.47
1:B:160:LYS:C	1:B:162:ALA:H	2.22	0.47
1:B:69:PRO:CG	4:B:202:HOH:O	2.41	0.47
1:B:5:ILE:CA	4:B:241:HOH:O	2.63	0.47
1:B:66:THR:CG2	4:B:174:HOH:O	2.63	0.47
1:B:90:GLU:C	1:B:92:ILE:N	2.71	0.47
1:B:112:TYR:OH	1:B:146:PHE:HA	2.15	0.47
1:B:2:ALA:CB	1:B:33:THR:HG21	2.45	0.47
1:B:32:LEU:CD1	1:B:33:THR:O	2.63	0.47
1:B:73:PRO:O	1:B:78:GLN:CG	2.56	0.47
1:B:93:SER:O	1:B:97:ARG:N	2.47	0.47
1:B:127:GLN:CA	1:B:127:GLN:OE1	2.63	0.47
1:B:146:PHE:O	1:B:148:ILE:HD11	2.15	0.47
1:A:59:GLU:O	1:A:135:ASN:ND2	2.47	0.47
1:A:90:GLU:C	1:A:92:ILE:N	2.71	0.47
1:A:68:VAL:HG13	1:A:69:PRO:HD2	1.97	0.46
1:B:75:PRO:N	1:B:78:GLN:CG	2.78	0.46
1:B:67:VAL:HG22	1:B:103:GLN:CB	2.34	0.46
1:B:69:PRO:CD	4:B:202:HOH:O	2.62	0.46
1:B:112:TYR:OH	1:B:147:TRP:N	2.48	0.46
1:B:112:TYR:CE1	1:B:146:PHE:CD1	3.04	0.46
1:A:147:TRP:NE1	1:A:160:LYS:HA	2.30	0.46
1:B:34:VAL:O	1:B:58:VAL:CG2	2.63	0.46
1:B:113:VAL:HB	4:B:244:HOH:O	2.14	0.46
1:B:141:LYS:HB3	1:B:164:GLY:O	2.16	0.46
1:B:43:ASP:O	1:B:47:GLY:N	2.49	0.46
1:B:64:VAL:CG1	1:B:128:LEU:CD1	2.94	0.46
1:B:79:ILE:O	1:B:79:ILE:CG2	2.61	0.46
1:B:129:LYS:H	1:B:129:LYS:HG2	1.20	0.46
1:B:10:PHE:HB2	1:B:30:PHE:CE1	2.51	0.46
1:B:142:TYR:CZ	1:B:144:PRO:CG	2.99	0.46
1:A:157:GLN:OE1	1:A:162:ALA:HB1	2.16	0.45
1:A:67:VAL:HG22	1:A:103:GLN:HB3	1.98	0.45
1:B:74:PRO:CB	1:B:75:PRO:HD2	2.46	0.45
1:B:147:TRP:O	1:B:148:ILE:HG13	2.15	0.45
1:B:90:GLU:CB	1:B:93:SER:HB3	2.46	0.45
1:B:105:VAL:HA	1:B:109:GLY:O	2.17	0.45
1:A:147:TRP:CD1	1:A:160:LYS:HA	2.51	0.45
1:B:75:PRO:C	1:B:77:ARG:N	2.75	0.45
1:B:154:CYS:SG	1:B:155:CYS:SG	3.14	0.45
1:B:159:ALA:O	1:B:162:ALA:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:169:GLU:O	1:B:170:ASN:ND2	2.49	0.45
1:B:99:PRO:CD	1:B:100:TYR:CD2	2.97	0.45
1:B:58:VAL:C	1:B:61:ILE:HG13	2.41	0.45
1:B:142:TYR:CZ	1:B:144:PRO:HB3	2.52	0.45
1:B:154:CYS:HB2	2:B:171:ZN:ZN	1.48	0.45
1:A:94:ILE:CD1	1:B:91:GLN:HG3	2.47	0.45
1:B:43:ASP:HB3	1:B:48:ARG:HG3	1.98	0.45
1:B:75:PRO:HB2	1:B:76:GLU:H	1.26	0.45
1:B:112:TYR:CE1	1:B:146:PHE:CE1	3.05	0.45
1:B:146:PHE:HB3	1:B:148:ILE:HG12	1.98	0.45
1:B:166:GLN:CG	1:B:167:ILE:N	2.78	0.45
1:B:122:LYS:N	4:B:192:HOH:O	2.46	0.44
1:B:143:HIS:CE1	1:B:154:CYS:HG	2.35	0.44
1:A:36:LYS:HA	1:A:58:VAL:HG23	1.98	0.44
1:A:91:GLN:HG3	1:B:94:ILE:HD13	1.99	0.44
1:B:133:ARG:NE	1:B:134:TYR:CE1	2.86	0.44
1:A:117:THR:HB	4:A:264:HOH:O	2.16	0.44
1:B:98:PHE:C	1:B:115:SER:O	2.61	0.44
1:B:142:TYR:CE1	1:B:144:PRO:CB	3.01	0.44
1:B:143:HIS:C	1:B:145:CYS:H	2.25	0.44
1:A:133:ARG:HB3	1:A:134:TYR:CD2	2.52	0.44
1:A:142:TYR:OH	1:A:144:PRO:HB3	2.17	0.44
1:B:11:LEU:HD12	1:B:11:LEU:HA	1.60	0.44
1:A:128:LEU:O	1:A:129:LYS:C	2.60	0.44
1:B:74:PRO:C	1:B:78:GLN:CG	2.91	0.44
1:B:147:TRP:C	1:B:147:TRP:CE3	2.96	0.44
1:A:90:GLU:O	1:A:93:SER:N	2.50	0.44
1:B:68:VAL:O	1:B:68:VAL:CG1	2.65	0.44
1:B:112:TYR:CD1	1:B:146:PHE:CE1	3.06	0.44
1:B:150:GLY:O	1:B:158:THR:CG2	2.66	0.44
1:A:89:MET:HE3	1:A:89:MET:O	2.18	0.44
1:B:13:ARG:HG2	4:B:184:HOH:O	2.17	0.43
1:B:36:LYS:CA	1:B:58:VAL:HG23	2.48	0.43
1:B:98:PHE:HA	1:B:115:SER:C	2.43	0.43
1:A:79:ILE:H	1:A:79:ILE:HG13	1.64	0.43
1:B:128:LEU:O	1:B:129:LYS:C	2.61	0.43
1:B:169:GLU:CG	1:B:170:ASN:N	2.76	0.43
1:B:146:PHE:CB	1:B:148:ILE:HD11	2.48	0.43
1:A:89:MET:HE3	1:A:89:MET:HB2	1.83	0.43
1:B:100:TYR:CB	1:B:115:SER:HB3	2.28	0.43
1:A:10:PHE:HB3	1:A:114:PHE:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:PHE:CD2	1:A:44:PHE:C	2.94	0.43
1:B:34:VAL:O	1:B:58:VAL:HB	2.18	0.43
1:B:74:PRO:HG3	1:B:77:ARG:NH2	2.34	0.43
1:A:143:HIS:NE2	1:A:164:GLY:HA2	2.33	0.43
1:A:5:ILE:CD1	1:A:33:THR:CA	2.96	0.43
1:A:155:CYS:O	1:A:157:GLN:N	2.51	0.43
1:B:73:PRO:C	1:B:78:GLN:HG2	2.43	0.43
1:A:141:LYS:CE	1:A:163:MET:CE	2.97	0.43
1:B:65:GLU:O	1:B:103:GLN:N	2.51	0.43
1:A:118:GLU:HG2	4:A:257:HOH:O	2.19	0.43
1:B:68:VAL:HA	1:B:69:PRO:HD2	1.79	0.43
1:B:64:VAL:CG1	1:B:128:LEU:HB3	2.46	0.43
1:B:75:PRO:N	1:B:78:GLN:CD	2.77	0.43
1:B:132:ILE:O	1:B:135:ASN:N	2.52	0.43
1:B:142:TYR:O	1:B:144:PRO:HD3	2.19	0.43
1:B:49:ARG:HG2	1:B:49:ARG:NH1	2.32	0.42
1:B:139:VAL:H	1:B:139:VAL:HG13	1.42	0.42
1:A:119:GLU:O	1:A:123:ARG:HG3	2.19	0.42
1:B:66:THR:HA	4:B:174:HOH:O	2.19	0.42
1:B:101:PRO:HB3	1:B:114:PHE:CE2	2.54	0.42
1:B:125:ILE:CG2	1:B:126:HIS:N	2.53	0.42
1:B:147:TRP:HA	1:B:152:TYR:HA	2.01	0.42
1:A:169:GLU:C	1:A:170:ASN:ND2	2.77	0.42
1:A:170:ASN:HD22	1:A:170:ASN:N	2.16	0.42
1:B:143:HIS:NE2	1:B:164:GLY:HA2	2.33	0.42
1:B:75:PRO:HA	1:B:78:GLN:CG	2.49	0.42
1:A:75:PRO:HA	1:A:78:GLN:HB2	2.02	0.42
1:B:11:LEU:HD12	1:B:26:LYS:O	2.20	0.42
1:B:143:HIS:CE1	1:B:163:MET:O	2.73	0.42
1:B:128:LEU:HD23	1:B:128:LEU:HA	1.74	0.42
1:A:25:PHE:C	1:A:26:LYS:HG2	2.45	0.42
1:A:92:ILE:HD13	1:A:92:ILE:HA	1.81	0.42
1:B:75:PRO:CA	1:B:78:GLN:CD	2.93	0.42
1:A:141:LYS:HA	1:A:166:GLN:O	2.20	0.41
1:A:145:CYS:HA	4:A:191:HOH:O	2.20	0.41
1:B:101:PRO:C	4:B:194:HOH:O	2.62	0.41
1:B:116:PRO:C	1:B:117:THR:CG2	2.92	0.41
1:A:42:TYR:CD2	1:B:44:PHE:CE2	3.08	0.41
1:B:17:LYS:HE3	1:B:106:TYR:OH	2.19	0.41
1:B:155:CYS:HG	1:B:165:CYS:HB3	1.75	0.41
1:B:4:VAL:HG13	1:B:6:LEU:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:THR:O	1:B:158:THR:HG22	2.20	0.41
1:B:160:LYS:C	1:B:162:ALA:N	2.78	0.41
1:B:26:LYS:HD2	1:B:28:ARG:NH2	2.35	0.41
1:B:106:TYR:CE2	1:B:108:GLU:HB2	2.54	0.41
1:A:9:ILE:HD12	1:B:92:ILE:HG12	2.02	0.41
1:A:74:PRO:HB2	1:A:75:PRO:HD2	2.03	0.41
1:B:142:TYR:CE1	1:B:144:PRO:HB3	2.55	0.41
1:A:126:HIS:O	1:A:127:GLN:C	2.61	0.41
1:A:155:CYS:C	1:A:157:GLN:N	2.79	0.41
1:B:98:PHE:HD2	1:B:116:PRO:HD3	1.86	0.41
1:B:122:LYS:O	1:B:123:ARG:C	2.61	0.41
1:B:114:PHE:CD1	4:B:178:HOH:O	2.58	0.40
1:B:143:HIS:HB3	1:B:154:CYS:SG	2.60	0.40
1:B:147:TRP:CH2	1:B:150:GLY:O	2.75	0.40
1:B:13:ARG:NH2	1:B:112:TYR:CD2	2.85	0.40
1:B:16:GLN:O	1:B:17:LYS:C	2.64	0.40
1:A:27:LYS:C	1:A:28:ARG:HG2	2.46	0.40
1:A:40:TYR:CD1	1:A:49:ARG:HD2	2.57	0.40
1:A:90:GLU:CG	1:A:91:GLN:N	2.84	0.40
1:A:122:LYS:O	1:A:123:ARG:C	2.62	0.40
1:B:25:PHE:CZ	1:B:77:ARG:CZ	3.04	0.40
1:B:87:SER:OG	1:B:117:THR:OG1	2.22	0.40
1:B:100:TYR:N	1:B:100:TYR:HD2	2.17	0.40
1:B:129:LYS:C	1:B:131:VAL:N	2.80	0.40
1:A:90:GLU:HA	1:A:93:SER:HG	1.87	0.40
1:B:66:THR:HG22	4:B:174:HOH:O	2.21	0.40
1:B:88:GLU:HB2	1:B:117:THR:HG22	1.99	0.40

All (7) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:207:HOH:O	4:A:207:HOH:O[6_555]	1.28	0.92
4:A:211:HOH:O	4:A:211:HOH:O[6_555]	1.70	0.50
1:A:142:TYR:OH	4:A:211:HOH:O[6_555]	1.76	0.44
1:A:129:LYS:NZ	1:A:170:ASN:O[6_555]	1.80	0.40
4:A:288:HOH:O	4:A:288:HOH:O[7_455]	1.91	0.29
1:A:66:THR:OG1	1:A:166:GLN:NE2[6_555]	2.07	0.13
4:B:201:HOH:O	4:B:201:HOH:O[10_555]	2.11	0.09

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	159/169 (94%)	143 (90%)	10 (6%)	6 (4%)	2	2
1	B	158/169 (94%)	103 (65%)	33 (21%)	22 (14%)	0	0
All	All	317/338 (94%)	246 (78%)	43 (14%)	28 (9%)	0	0

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	B	75	PRO
1	B	76	GLU
1	B	78	GLN
1	B	91	GLN
1	B	125	ILE
1	B	133	ARG
1	B	157	GLN
1	B	3	ALA
1	B	17	LYS
1	B	110	PRO
1	B	130	ASN
1	B	152	TYR
1	B	156	SER
1	B	166	GLN
1	A	160	LYS
1	B	7	GLU
1	B	69	PRO
1	B	144	PRO
1	A	110	PRO
1	A	133	ARG
1	B	126	HIS
1	B	153	LEU
1	B	118	GLU
1	A	138	LEU

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Mol	Chain	Res	Type
1	B	139	VAL
1	B	100	TYR
1	A	80	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	153/158 (97%)	116 (76%)	37 (24%)	1	1
1	B	152/158 (96%)	106 (70%)	46 (30%)	0	0
All	All	305/316 (96%)	222 (73%)	83 (27%)	0	0

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ILE
1	A	12	LYS
1	A	17	LYS
1	A	18	LYS
1	A	19	LYS
1	A	20	THR
1	A	25	PHE
1	A	36	LYS
1	A	38	SER
1	A	44	PHE
1	A	49	ARG
1	A	51	SER
1	A	68	VAL
1	A	70	GLU
1	A	77	ARG
1	A	79	ILE
1	A	81	ARG
1	A	82	ARG
1	A	89	MET
1	A	90	GLU

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Mol	Chain	Res	Type
1	A	94	ILE
1	A	96	GLU
1	A	97	ARG
1	A	105	VAL
1	A	107	ASP
1	A	118	GLU
1	A	119	GLU
1	A	120	LEU
1	A	122	LYS
1	A	133	ARG
1	A	139	VAL
1	A	140	GLN
1	A	148	ILE
1	A	151	GLN
1	A	160	LYS
1	A	169	GLU
1	A	170	ASN
1	B	5	ILE
1	B	6	LEU
1	B	12	LYS
1	B	15	GLN
1	B	16	GLN
1	B	17	LYS
1	B	25	PHE
1	B	32	LEU
1	B	34	VAL
1	B	36	LYS
1	B	37	LEU
1	B	38	SER
1	B	40	TYR
1	B	59	GLU
1	B	60	LYS
1	B	61	ILE
1	B	62	THR
1	B	65	GLU
1	B	68	VAL
1	B	70	GLU
1	B	94	ILE
1	B	105	VAL
1	B	108	GLU
1	B	114	PHE
1	B	117	THR

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Mol	Chain	Res	Type
1	B	119	GLU
1	B	120	LEU
1	B	125	ILE
1	B	126	HIS
1	B	131	VAL
1	B	133	ARG
1	B	136	SER
1	B	137	ASP
1	B	139	VAL
1	B	140	GLN
1	B	145	CYS
1	B	148	ILE
1	B	153	LEU
1	B	155	CYS
1	B	157	GLN
1	B	158	THR
1	B	160	LYS
1	B	163	MET
1	B	165	CYS
1	B	168	LEU
1	B	169	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	103	GLN
1	A	135	ASN
1	A	170	ASN
1	B	35	HIS
1	B	126	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	4IP	A	171	-	28,28,28	2.13	13 (46%)	46,46,46	1.72	9 (19%)
3	4IP	B	172	-	28,28,28	2.09	15 (53%)	46,46,46	2.19	18 (39%)

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	4IP	A	171	-	-	0/20/44/44	0/1/1/1
3	4IP	B	172	-	-	3/20/44/44	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	171	4IP	P4-O4	-4.58	1.51	1.59
3	B	172	4IP	P1-O1	3.84	1.66	1.59
3	B	172	4IP	P4-O4	-3.79	1.52	1.59
3	A	171	4IP	P1-O1	3.72	1.66	1.59
3	A	171	4IP	P3-O4P	-3.10	1.40	1.50
3	A	171	4IP	P3-O6P	-3.08	1.43	1.54
3	B	172	4IP	P5-OPH	-2.96	1.43	1.54
3	B	172	4IP	P4-O7P	-2.90	1.44	1.54
3	A	171	4IP	P5-OPH	-2.86	1.44	1.54
3	A	171	4IP	P3-O3	-2.82	1.54	1.59
3	B	172	4IP	P3-O5P	-2.81	1.44	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	172	4IP	O3-C3	-2.60	1.35	1.44
3	A	171	4IP	P3-O5P	-2.56	1.45	1.54
3	B	172	4IP	P3-O6P	-2.48	1.45	1.54
3	B	172	4IP	O1-C1	2.42	1.52	1.44
3	A	171	4IP	C5-C4	-2.40	1.47	1.52
3	B	172	4IP	C2-C3	-2.33	1.45	1.52
3	B	172	4IP	C4-C3	2.32	1.57	1.52
3	B	172	4IP	P3-O4P	-2.23	1.43	1.50
3	A	171	4IP	C4-C3	2.21	1.56	1.52
3	A	171	4IP	P5-OPF	-2.15	1.46	1.54
3	A	171	4IP	P4-O9P	-2.11	1.43	1.50
3	A	171	4IP	O1-C1	2.11	1.51	1.44
3	B	172	4IP	P1-O3P	-2.10	1.47	1.54
3	B	172	4IP	P4-O9P	2.10	1.57	1.50
3	B	172	4IP	P3-O3	-2.08	1.55	1.59
3	B	172	4IP	P5-OPG	-2.05	1.44	1.50
3	A	171	4IP	P5-OPG	-2.02	1.44	1.50

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	172	4IP	O3-P3-O4P	-5.52	89.64	109.33
3	B	172	4IP	O1-C1-C2	5.29	119.92	108.73
3	A	171	4IP	O1-C1-C2	5.07	119.47	108.73
3	B	172	4IP	O5P-P3-O4P	3.98	126.33	110.83
3	B	172	4IP	O7P-P4-O4	-3.89	90.69	105.85
3	B	172	4IP	O3-C3-C4	3.15	115.46	108.76
3	A	171	4IP	O3P-P1-O1	-3.09	93.82	105.85
3	B	172	4IP	OPH-P5-O5	-3.08	93.84	105.85
3	B	172	4IP	C6-C1-C2	-2.93	106.79	110.86
3	A	171	4IP	O3-P3-O4P	-2.93	98.91	109.33
3	B	172	4IP	O2-C2-C1	2.87	117.30	109.94
3	B	172	4IP	P5-O5-C5	2.87	131.09	123.43
3	A	171	4IP	C6-C1-C2	2.86	114.83	110.86
3	B	172	4IP	O5P-P3-O3	-2.85	94.75	105.85
3	B	172	4IP	OPH-P5-OPF	2.84	118.47	107.80
3	A	171	4IP	C3-C2-C1	-2.78	103.47	109.11
3	A	171	4IP	O5P-P3-O4P	2.72	121.45	110.83
3	A	171	4IP	P5-O5-C5	2.68	130.60	123.43
3	B	172	4IP	C5-C4-C3	-2.68	104.55	110.43
3	A	171	4IP	O3-C3-C4	2.68	114.45	108.76
3	B	172	4IP	P1-O1-C1	2.47	130.02	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	172	4IP	O5-C5-C4	2.37	113.81	108.76
3	A	171	4IP	O3P-P1-O1P	2.07	118.91	110.83
3	B	172	4IP	O1-P1-O1P	-2.05	102.04	109.33
3	B	172	4IP	O4-C4-C3	-2.04	104.41	108.76
3	B	172	4IP	O7P-P4-O9P	2.03	118.75	110.83
3	B	172	4IP	O2P-P1-O1P	2.03	118.75	110.83

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	172	4IP	C4-C5-O5-P5
3	B	172	4IP	C1-O1-P1-O3P
3	B	172	4IP	C1-O1-P1-O2P

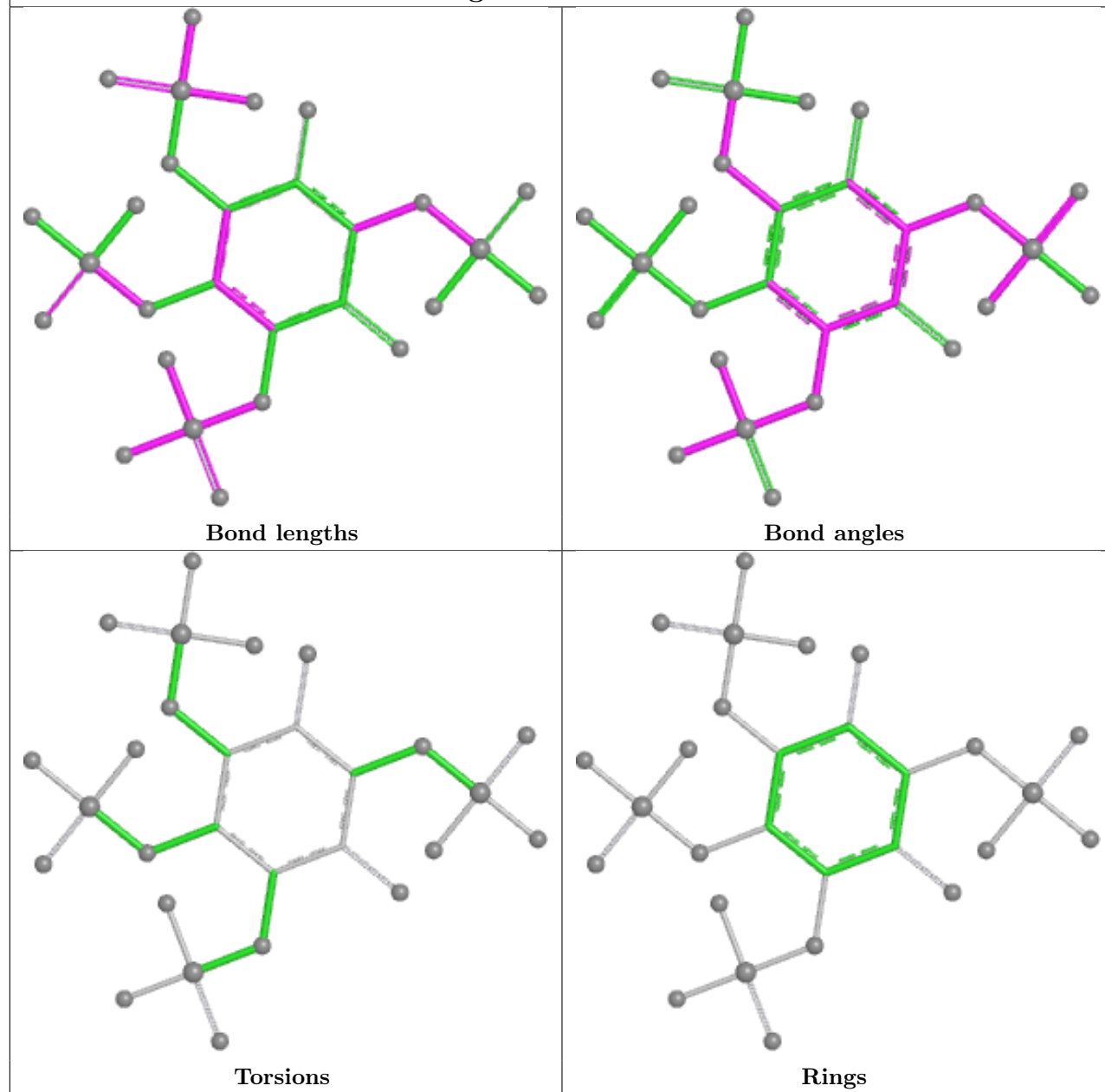
There are no ring outliers.

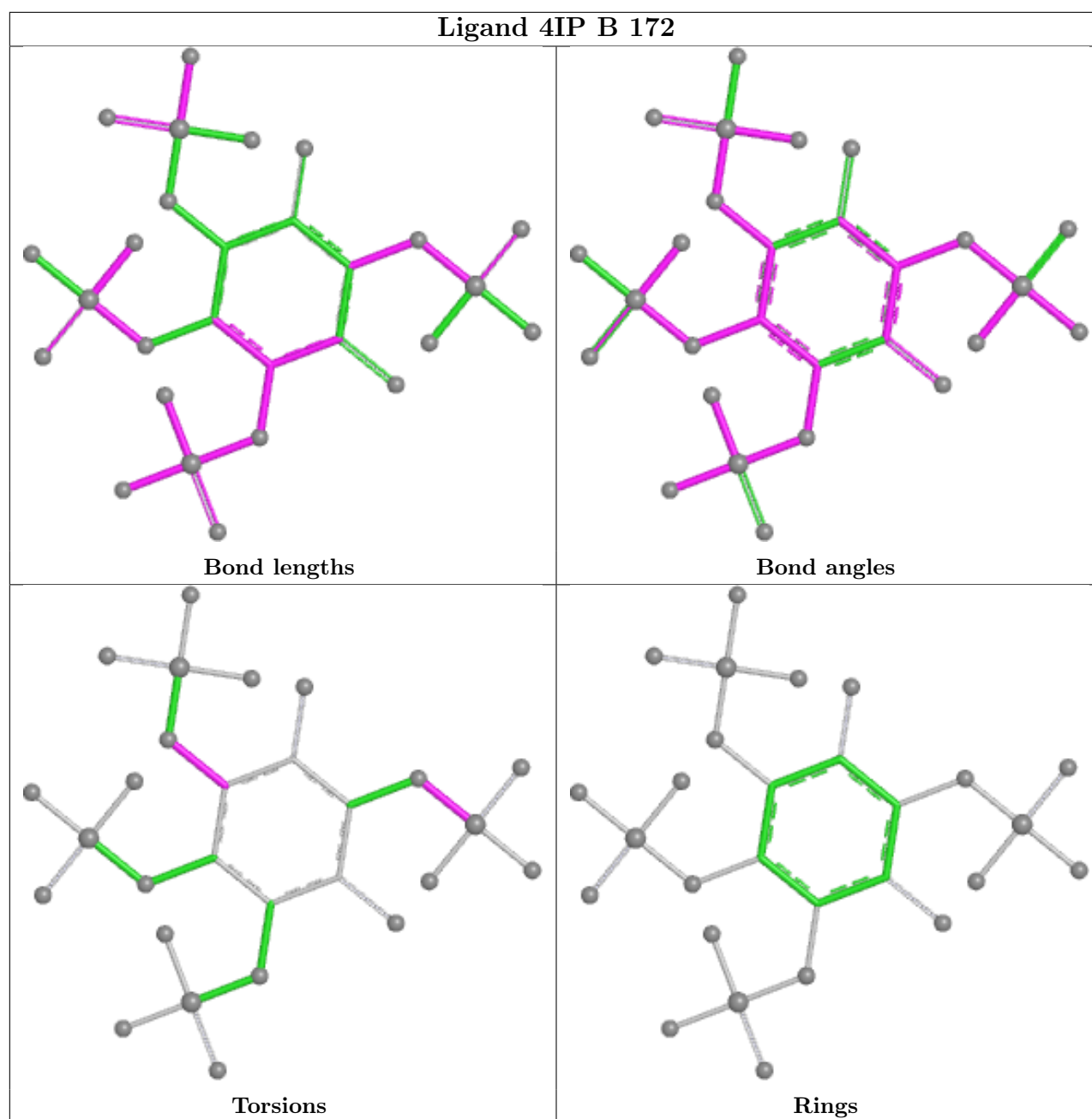
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	171	4IP	2	0
3	B	172	4IP	2	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 4IP A 171





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	163/169 (96%)	0.11	2 (1%) 76 73	33, 53, 77, 85	8 (4%)
1	B	162/169 (95%)	0.80	16 (9%) 13 9	16, 59, 82, 85	16 (9%)
All	All	325/338 (96%)	0.46	18 (5%) 30 27	16, 56, 80, 85	24 (7%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	TYR	4.9
1	B	155	CYS	3.8
1	B	87	SER	3.5
1	B	152	TYR	3.1
1	A	170	ASN	2.8
1	B	170	ASN	2.7
1	B	100	TYR	2.7
1	B	153	LEU	2.6
1	B	103	GLN	2.6
1	B	141	LYS	2.6
1	B	88	GLU	2.5
1	B	164	GLY	2.5
1	B	131	VAL	2.4
1	B	158	THR	2.3
1	A	89	MET	2.1
1	B	162	ALA	2.1
1	B	167	ILE	2.1
1	B	169	GLU	2.1

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](#)

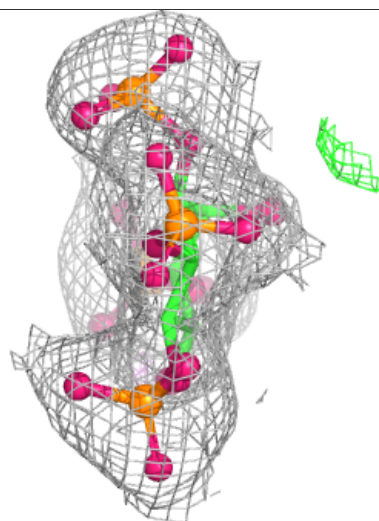
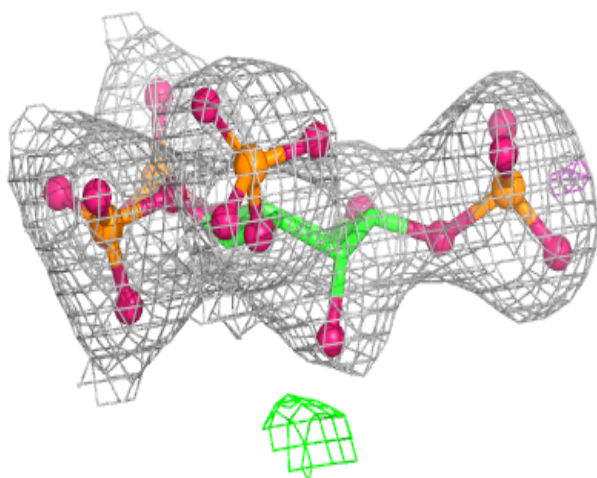
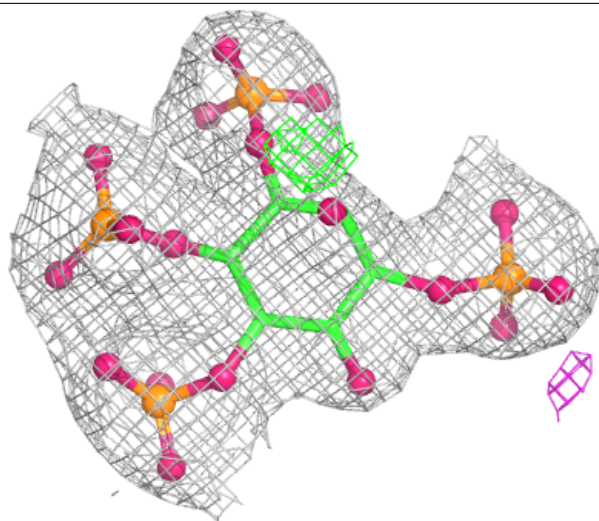
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ZN	B	171	1/1	0.93	0.13	85,85,85,85	0
3	4IP	B	172	28/28	0.97	0.07	34,56,63,75	0
3	4IP	A	171	28/28	0.98	0.05	35,44,54,65	0
2	ZN	A	1	1/1	1.00	0.02	45,45,45,45	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.

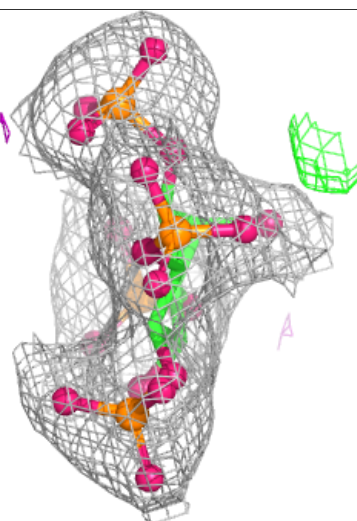
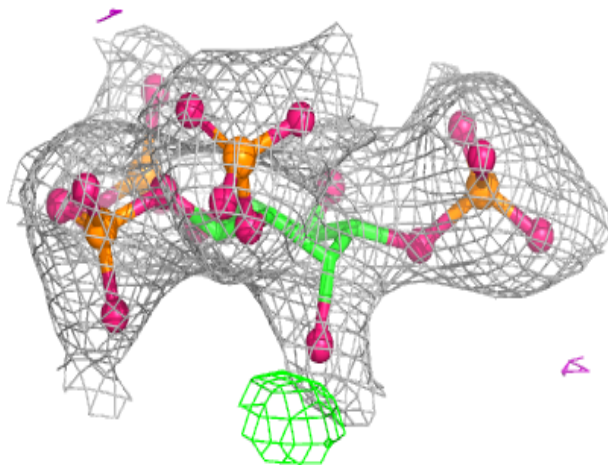
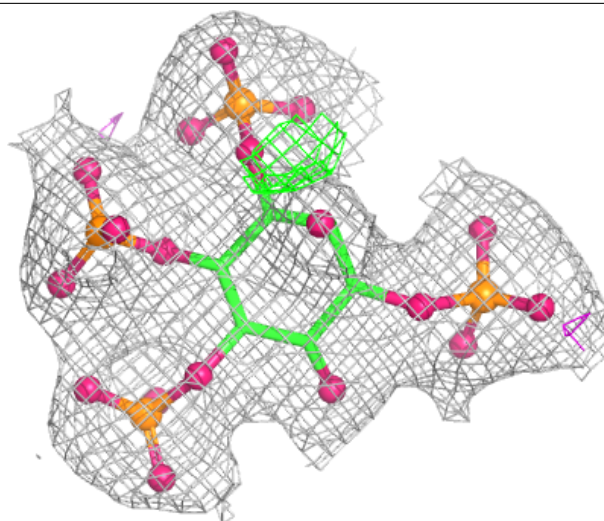
Electron density around 4IP B 172:

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



Electron density around 4IP A 171:

$2mF_o-DF_c$ (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.