



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:01 PM UTC

PDB ID : 2Z0P / pdb_00002z0p
Title : Crystal structure of PH domain of Bruton's tyrosine kinase
Authors : Murayama, K.; Kato-Murayama, M.; Mishima, C.; Shirouzu, M.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-05-07
Resolution : 2.58 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

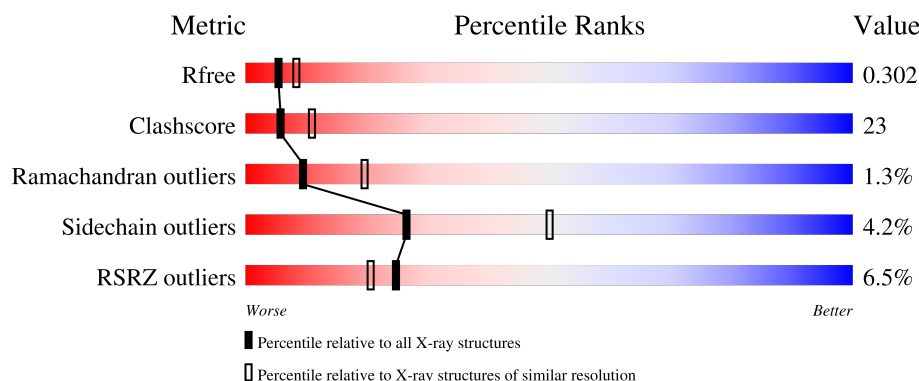
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

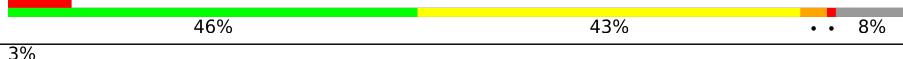
The reported resolution of this entry is 2.58 Å.

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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	
1	B	169	
1	C	169	
1	D	169	

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

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	4PT	B	502	X	-	-	-
3	4PT	D	504	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5510 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.

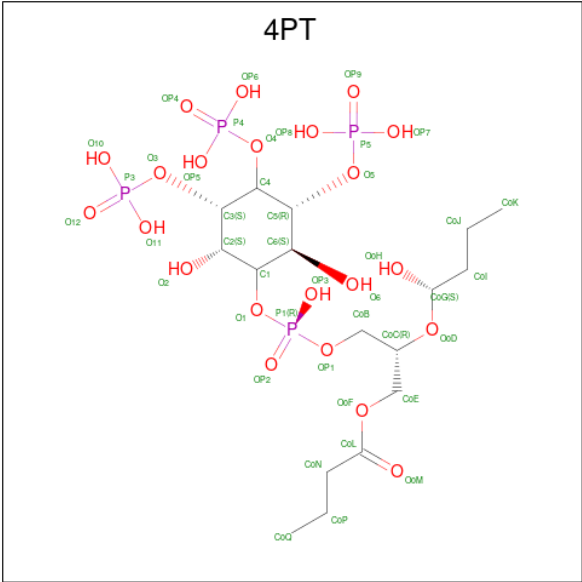
- Molecule 1 is a protein called Tyrosine-protein kinase BTK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	158	Total	C	N	O	S	0	0	0
			1316	851	226	232	7			
1	B	161	Total	C	N	O	S	0	0	0
			1343	865	230	241	7			
1	C	155	Total	C	N	O	S	0	0	0
			1296	836	222	231	7			
1	D	161	Total	C	N	O	S	0	0	0
			1341	865	230	239	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		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (2R)-3-{[(S)-{[(2S,3R,5S,6S)-2,6-DIHYDROXY-3,4,5-TRIS(PHOSPHONOXY)CYCLOHEXYL]OXY} (HYDROXY)PHOSPHORYL]OXY}-2-(1-HYDROXY BUTOXY)PROPYL BUTYRATE (CCD ID: 4PT) (formula: C₁₇H₃₆O₂₂P₄).



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
			43	17	22	4		
3	C	1	Total	C	O	P	0	0
			28	6	18	4		
3	D	1	Total	C	O	P	0	0
			43	17	22	4		

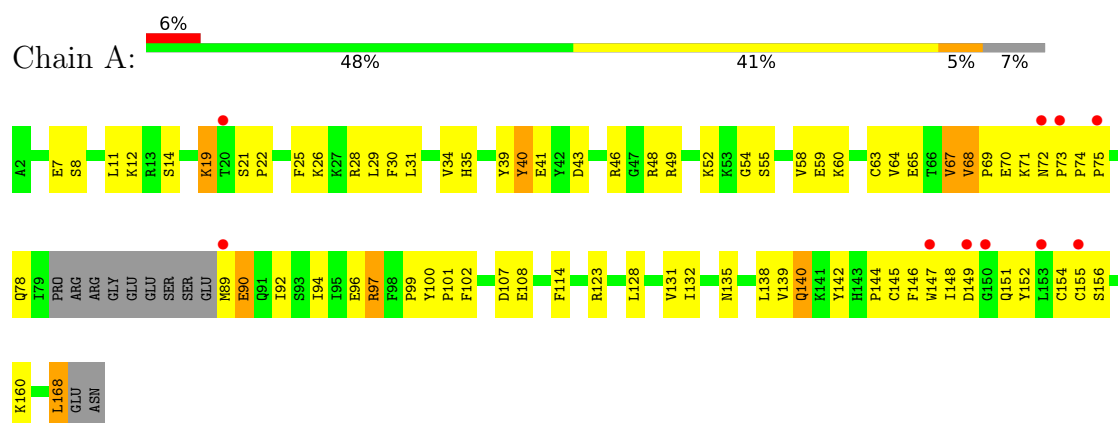
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total	O	0	0
			13	13		
4	B	25	Total	O	0	0
			25	25		
4	C	6	Total	O	0	0
			6	6		
4	D	24	Total	O	0	0
			24	24		

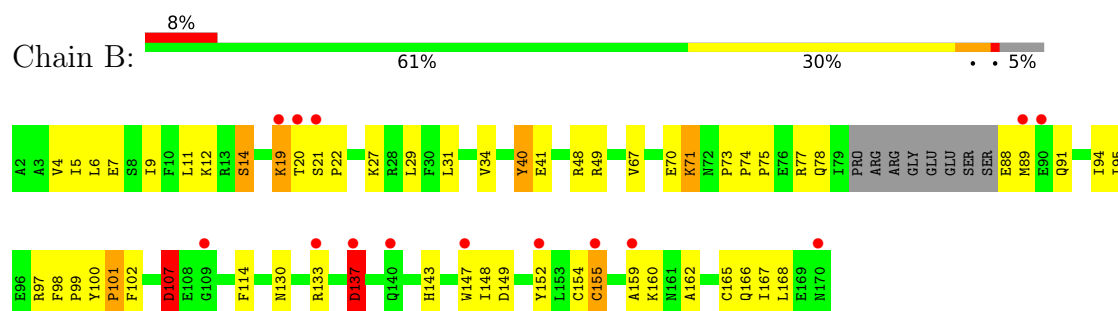
3 Residue-property plots [i](#)

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

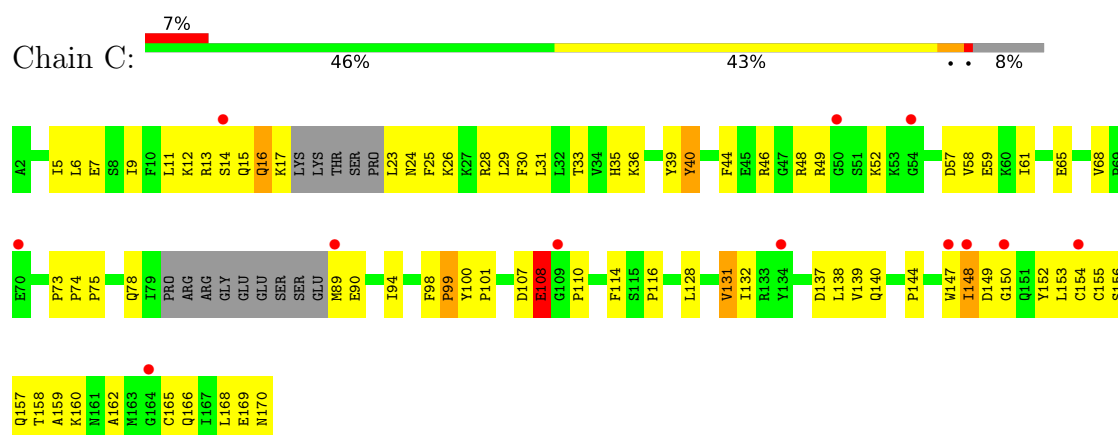
• Molecule 1: Tyrosine-protein kinase BTK



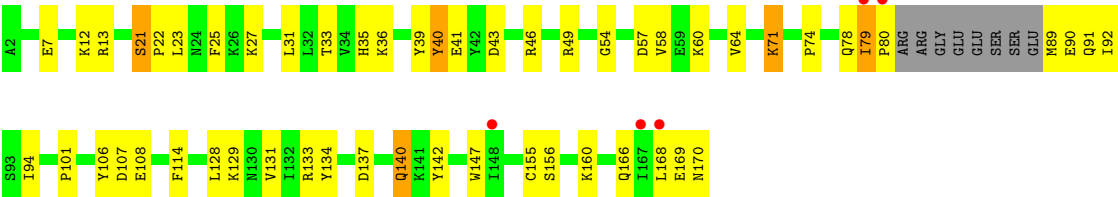
• Molecule 1: Tyrosine-protein kinase BTK



• Molecule 1: Tyrosine-protein kinase BTK



● Molecule 1: Tyrosine-protein kinase BTK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.42Å 62.50Å 117.06Å 90.00° 94.33° 90.00°	Depositor
Resolution (Å)	29.67 – 2.58 29.67 – 2.58	Depositor EDS
% Data completeness (in resolution range)	98.1 (29.67-2.58) 98.1 (29.67-2.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.84 (at 2.57Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.242 , 0.300 0.243 , 0.302	Depositor DCC
R_{free} test set	1283 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	44.1	Xtriage
Anisotropy	0.764	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5510	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 33.69 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.6864e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 4PT, 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	0.47	0/1348	0.98	4/1817 (0.2%)
1	B	0.47	0/1375	1.00	5/1852 (0.3%)
1	C	0.42	0/1326	0.92	2/1785 (0.1%)
1	D	0.50	0/1374	0.99	5/1852 (0.3%)
All	All	0.47	0/5423	0.97	16/7306 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	155	CYS	N-CA-C	-7.92	98.29	109.69
1	D	79	ILE	CB-CA-C	-7.51	106.57	113.70
1	D	41	GLU	N-CA-C	-6.84	100.16	110.28
1	A	90	GLU	N-CA-C	6.12	118.45	111.11
1	B	19	LYS	N-CA-C	5.96	123.50	110.80
1	B	41	GLU	N-CA-C	-5.80	101.70	110.28
1	B	137	ASP	N-CA-C	5.78	120.11	112.25
1	A	68	VAL	CA-C-N	5.58	125.53	119.78
1	A	68	VAL	C-N-CA	5.58	125.53	119.78
1	A	41	GLU	N-CA-C	-5.41	101.66	110.20
1	C	131	VAL	N-CA-C	5.21	117.61	111.09
1	D	21	SER	CA-C-N	5.20	125.65	120.14
1	D	21	SER	C-N-CA	5.20	125.65	120.14
1	D	33	THR	N-CA-C	-5.17	102.54	110.14
1	C	166	GLN	N-CA-C	5.16	117.23	109.23
1	B	4	VAL	N-CA-C	5.00	114.81	107.15

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	1316	0	1333	81	0
1	B	1343	0	1352	55	0
1	C	1296	0	1299	71	0
1	D	1341	0	1352	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	28	0	8	1	0
3	B	43	0	27	5	0
3	C	28	0	8	1	0
3	D	43	0	27	4	0
4	A	13	0	0	0	0
4	B	25	0	0	3	0
4	C	6	0	0	0	0
4	D	24	0	0	1	0
All	All	5510	0	5406	252	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 23.

All (252) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:VAL:HG11	1:A:131:VAL:HG12	1.29	1.12
1:B:71:LYS:HE3	1:B:71:LYS:H	0.99	1.10
1:C:159:ALA:HB3	1:C:162:ALA:HB2	1.33	1.10
1:B:159:ALA:HB3	1:B:162:ALA:HB2	1.34	1.05
1:B:71:LYS:HE3	1:B:71:LYS:N	1.71	1.05
1:A:97:ARG:HH11	1:A:97:ARG:HB2	1.20	1.04
1:D:71:LYS:HE3	1:D:71:LYS:H	1.22	1.00
1:A:89:MET:HE3	1:A:90:GLU:H	1.27	0.96
1:C:148:ILE:HD12	1:C:153:LEU:HD21	1.48	0.96
1:A:19:LYS:H	1:A:19:LYS:HD2	1.30	0.96
1:D:140:GLN:H	1:D:140:GLN:CD	1.78	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:58:VAL:HG11	1:D:131:VAL:HG12	1.53	0.90
1:A:58:VAL:HG11	1:A:131:VAL:CG1	2.01	0.89
1:C:58:VAL:HG11	1:C:131:VAL:HG12	1.52	0.89
1:A:26:LYS:HD3	1:A:28:ARG:HH22	1.42	0.83
1:D:71:LYS:H	1:D:71:LYS:CE	1.92	0.81
1:A:123:ARG:HH12	1:B:89:MET:HE2	1.47	0.80
1:B:130:ASN:O	1:B:133:ARG:HG2	1.85	0.77
1:A:19:LYS:H	1:A:19:LYS:CD	1.94	0.76
1:D:58:VAL:HG11	1:D:131:VAL:CG1	2.16	0.75
1:C:152:TYR:HB3	1:C:154:CYS:SG	2.25	0.75
1:B:29:LEU:HD23	1:B:49:ARG:HG3	1.67	0.75
1:D:89:MET:HE3	1:D:90:GLU:HB2	1.68	0.74
1:A:74:PRO:O	1:A:78:GLN:HG3	1.88	0.72
1:D:71:LYS:HE3	1:D:71:LYS:N	2.01	0.72
1:C:40:TYR:CE2	1:C:52:LYS:HG2	2.23	0.72
3:D:504:4PT:H0N2	3:D:504:4PT:H0C	1.72	0.71
1:C:17:LYS:HE2	1:C:108:GLU:HG2	1.72	0.71
1:C:89:MET:HG2	1:C:90:GLU:H	1.55	0.71
1:A:97:ARG:HB2	1:A:97:ARG:NH1	2.00	0.70
1:C:58:VAL:CG1	1:C:131:VAL:HG12	2.22	0.69
1:B:74:PRO:O	1:B:78:GLN:HG3	1.93	0.69
3:D:504:4PT:H0E1	3:D:504:4PT:O0H	1.92	0.68
1:A:58:VAL:CG1	1:A:131:VAL:HG12	2.17	0.68
1:C:11:LEU:HB3	1:C:114:PHE:HB2	1.76	0.67
1:D:36:LYS:HA	1:D:58:VAL:HG23	1.76	0.66
1:C:12:LYS:HE2	3:C:503:4PT:O10	1.95	0.66
1:A:26:LYS:HB3	1:A:28:ARG:NH1	2.10	0.66
1:C:148:ILE:CD1	1:C:153:LEU:HD21	2.25	0.66
1:D:43:ASP:OD2	1:D:46:ARG:HB2	1.95	0.66
1:C:58:VAL:HG11	1:C:131:VAL:CG1	2.24	0.65
1:C:89:MET:HG2	1:C:90:GLU:N	2.11	0.64
1:B:7:GLU:HG3	1:B:31:LEU:CD2	2.27	0.64
1:B:75:PRO:N	1:B:78:GLN:HE21	1.96	0.64
3:B:502:4PT:O12	3:B:502:4PT:OP5	2.16	0.63
1:A:92:ILE:HD11	1:B:9:ILE:HG13	1.80	0.63
1:B:88:GLU:HB3	4:B:520:HOH:O	1.99	0.63
1:C:101:PRO:HB3	1:C:114:PHE:CD2	2.32	0.63
1:A:26:LYS:HD3	1:A:28:ARG:NH2	2.14	0.62
1:A:123:ARG:NH1	1:B:89:MET:HE2	2.12	0.62
1:B:97:ARG:HD2	4:B:512:HOH:O	1.97	0.62
1:B:12:LYS:HE2	3:B:502:4PT:O10	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:PRO:O	1:C:101:PRO:HD3	2.00	0.62
1:D:7:GLU:HG3	1:D:31:LEU:CD2	2.29	0.62
1:A:26:LYS:HB3	1:A:28:ARG:HH12	1.64	0.61
1:C:101:PRO:HB3	1:C:114:PHE:CE2	2.36	0.61
1:A:101:PRO:HB3	1:A:114:PHE:CE2	2.35	0.61
1:B:11:LEU:HB3	1:B:114:PHE:HB2	1.82	0.61
1:B:74:PRO:C	1:B:78:GLN:HE21	2.10	0.60
1:B:11:LEU:HD23	1:B:77:ARG:HD3	1.84	0.60
1:B:147:TRP:CE3	1:B:160:LYS:HE3	2.37	0.59
1:A:19:LYS:HD2	1:A:19:LYS:N	2.08	0.59
1:A:140:GLN:CD	1:A:140:GLN:H	2.10	0.58
1:A:97:ARG:HH11	1:A:97:ARG:CB	2.07	0.58
1:B:71:LYS:C	1:B:73:PRO:HD3	2.28	0.58
1:B:70:GLU:HB2	1:B:73:PRO:HG3	1.86	0.58
1:C:74:PRO:C	1:C:78:GLN:HE21	2.11	0.58
1:D:140:GLN:CD	1:D:140:GLN:N	2.56	0.58
1:A:140:GLN:O	1:A:168:LEU:HB2	2.04	0.58
1:D:79:ILE:HB	1:D:80:PRO:HD3	1.83	0.58
1:C:5:ILE:HG22	1:C:6:LEU:HG	1.85	0.57
1:C:15:GLN:O	1:C:16:GLN:HB2	2.04	0.57
1:A:46:ARG:HD3	1:A:48:ARG:NH2	2.18	0.57
1:A:142:TYR:CD1	1:A:168:LEU:HD22	2.39	0.57
1:B:130:ASN:HA	1:B:133:ARG:HE	1.69	0.57
1:C:57:ASP:HB3	1:C:59:GLU:HG2	1.86	0.57
1:C:148:ILE:HD11	1:C:153:LEU:HD11	1.86	0.57
1:A:43:ASP:OD2	1:A:46:ARG:HD2	2.04	0.56
1:B:148:ILE:HG22	1:B:149:ASP:OD2	2.05	0.56
1:A:97:ARG:HG2	1:A:99:PRO:HD3	1.86	0.56
1:A:147:TRP:HB2	1:A:152:TYR:CE1	2.39	0.56
1:A:34:VAL:HB	1:D:134:TYR:CD2	2.40	0.56
1:C:17:LYS:HE2	1:C:108:GLU:CG	2.35	0.55
1:B:154:CYS:HB2	1:B:165:CYS:SG	2.46	0.55
1:A:148:ILE:O	1:A:149:ASP:HB2	2.07	0.55
1:C:94:ILE:HG23	1:D:91:GLN:HG3	1.88	0.54
1:A:64:VAL:O	1:A:65:GLU:HG2	2.07	0.54
1:C:155:CYS:SG	1:C:157:GLN:HG3	2.48	0.54
1:A:75:PRO:N	1:A:78:GLN:HE21	2.05	0.53
1:D:142:TYR:CE1	1:D:168:LEU:HD23	2.43	0.53
1:C:58:VAL:HA	1:C:61:ILE:HD12	1.90	0.53
1:C:40:TYR:CD2	1:C:49:ARG:HD3	2.43	0.53
1:A:60:LYS:HE3	1:A:107:ASP:OD1	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:PRO:HG2	1:C:78:GLN:HG2	1.89	0.53
1:A:147:TRP:O	1:A:147:TRP:CE3	2.61	0.53
1:A:92:ILE:O	1:A:96:GLU:HG3	2.08	0.52
1:D:107:ASP:OD2	1:D:108:GLU:HG3	2.09	0.52
1:A:40:TYR:CD1	1:A:40:TYR:N	2.77	0.52
1:A:101:PRO:HB3	1:A:114:PHE:CD2	2.44	0.52
1:A:107:ASP:OD2	1:A:108:GLU:HG3	2.09	0.52
1:A:148:ILE:HG12	1:A:149:ASP:OD1	2.09	0.52
1:B:5:ILE:HG22	1:B:6:LEU:HG	1.91	0.52
1:B:155:CYS:HB3	1:B:165:CYS:CB	2.40	0.52
1:C:9:ILE:HG13	1:D:92:ILE:HD11	1.90	0.51
1:D:40:TYR:CD2	1:D:49:ARG:HD2	2.46	0.51
1:A:145:CYS:HB2	1:A:154:CYS:HB3	1.93	0.51
1:C:30:PHE:O	1:C:31:LEU:HD23	2.11	0.51
1:C:128:LEU:O	1:C:132:ILE:HG12	2.10	0.51
1:C:170:ASN:HD22	1:C:170:ASN:C	2.18	0.51
1:B:99:PRO:HG2	1:B:100:TYR:CE2	2.46	0.51
1:A:71:LYS:C	1:A:73:PRO:HD3	2.36	0.51
3:B:502:4PT:H0E2	3:B:502:4PT:H0P2	1.93	0.51
1:D:71:LYS:HE2	4:D:508:HOH:O	2.10	0.51
1:D:89:MET:CE	1:D:90:GLU:HB2	2.39	0.51
1:A:29:LEU:HD23	1:A:49:ARG:HB3	1.93	0.50
1:A:46:ARG:HD3	1:A:48:ARG:CZ	2.41	0.50
1:C:68:VAL:HG23	1:C:144:PRO:O	2.12	0.50
1:D:89:MET:HG3	1:D:90:GLU:N	2.27	0.50
1:C:44:PHE:O	1:D:27:LYS:HE3	2.12	0.50
1:C:147:TRP:NE1	1:C:160:LYS:HA	2.27	0.49
1:C:25:PHE:O	1:C:26:LYS:HG2	2.13	0.49
1:A:74:PRO:C	1:A:78:GLN:HE21	2.20	0.49
1:A:39:TYR:CZ	1:A:54:GLY:HA3	2.47	0.49
1:A:94:ILE:HD13	1:B:91:GLN:HB2	1.94	0.49
1:C:40:TYR:N	1:C:40:TYR:CD1	2.81	0.49
1:D:40:TYR:CD1	1:D:40:TYR:N	2.80	0.49
1:C:9:ILE:HG13	1:D:92:ILE:CD1	2.43	0.48
1:C:65:GLU:CD	1:C:168:LEU:HD11	2.38	0.48
1:B:101:PRO:HB3	1:B:114:PHE:CD2	2.48	0.48
1:C:7:GLU:HG3	1:C:31:LEU:CD2	2.43	0.48
1:D:12:LYS:HE3	1:D:39:TYR:CE2	2.48	0.48
1:A:59:GLU:HA	1:A:135:ASN:OD1	2.13	0.48
1:B:7:GLU:HG3	1:B:31:LEU:HD23	1.94	0.48
1:C:33:THR:HG1	1:C:36:LYS:H	1.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:504:4PT:H0N2	3:D:504:4PT:C0C	2.35	0.48
1:A:14:SER:OG	3:A:501:4PT:H4	2.14	0.47
1:B:167:ILE:HG12	1:B:168:LEU:N	2.29	0.47
1:B:107:ASP:OD2	1:B:107:ASP:N	2.47	0.47
1:A:12:LYS:NZ	1:A:39:TYR:CE2	2.76	0.47
1:D:107:ASP:OD2	1:D:108:GLU:N	2.48	0.47
1:C:169:GLU:O	1:C:170:ASN:HB2	2.14	0.47
1:D:101:PRO:HB3	1:D:114:PHE:CE2	2.49	0.47
1:A:8:SER:O	1:A:29:LEU:HD12	2.15	0.47
1:A:89:MET:HE3	1:A:90:GLU:N	2.11	0.47
1:C:152:TYR:O	1:C:156:SER:N	2.38	0.47
1:D:129:LYS:O	1:D:133:ARG:HG3	2.15	0.47
1:A:35:HIS:HA	1:D:35:HIS:HE1	1.79	0.47
1:A:67:VAL:HG12	1:A:102:PHE:HA	1.96	0.47
1:B:137:ASP:N	1:B:137:ASP:OD1	2.47	0.46
1:C:15:GLN:O	1:C:16:GLN:CB	2.63	0.46
1:B:34:VAL:HG22	4:B:508:HOH:O	2.15	0.46
1:B:101:PRO:HB3	1:B:114:PHE:CE2	2.51	0.46
1:A:25:PHE:CD1	1:A:25:PHE:N	2.83	0.46
1:B:71:LYS:O	1:B:73:PRO:HD3	2.14	0.46
1:D:89:MET:CG	1:D:90:GLU:H	2.28	0.46
1:C:29:LEU:O	1:C:39:TYR:HA	2.16	0.46
1:C:74:PRO:O	1:C:78:GLN:HG3	2.16	0.46
1:D:58:VAL:CG1	1:D:131:VAL:HG12	2.33	0.46
1:A:7:GLU:HG3	1:A:31:LEU:CD2	2.47	0.45
1:A:63:CYS:HB2	1:A:139:VAL:CG2	2.46	0.45
1:C:12:LYS:HG2	1:C:13:ARG:O	2.16	0.45
1:C:75:PRO:N	1:C:78:GLN:HE21	2.13	0.45
1:C:13:ARG:HG2	1:C:14:SER:N	2.32	0.45
1:A:75:PRO:HA	1:A:78:GLN:NE2	2.31	0.45
1:A:145:CYS:SG	1:A:154:CYS:HB3	2.57	0.45
1:A:155:CYS:O	1:A:156:SER:HB2	2.17	0.45
1:C:5:ILE:O	1:C:6:LEU:HD23	2.16	0.45
1:C:99:PRO:HG2	1:C:100:TYR:CE2	2.51	0.45
1:D:169:GLU:HA	1:D:169:GLU:OE2	2.16	0.45
1:A:52:LYS:HE3	1:A:55:SER:HB2	1.97	0.45
1:A:147:TRP:CZ3	1:A:160:LYS:NZ	2.85	0.45
1:A:147:TRP:O	1:A:147:TRP:HE3	2.00	0.45
1:C:138:LEU:O	1:C:139:VAL:C	2.60	0.45
1:D:89:MET:HG3	1:D:90:GLU:H	1.81	0.45
1:D:13:ARG:HD2	1:D:23:LEU:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:74:PRO:O	1:D:78:GLN:HG3	2.16	0.44
1:A:142:TYR:O	1:A:144:PRO:HD3	2.16	0.44
1:C:7:GLU:CG	1:C:29:LEU:HD11	2.47	0.44
1:D:21:SER:HA	1:D:22:PRO:HD3	1.76	0.44
1:C:46:ARG:HD2	1:C:48:ARG:NH2	2.32	0.44
1:C:154:CYS:HB2	1:C:165:CYS:SG	2.57	0.44
1:A:128:LEU:O	1:A:132:ILE:HG12	2.17	0.44
1:B:75:PRO:N	1:B:78:GLN:NE2	2.65	0.44
1:D:25:PHE:N	1:D:25:PHE:CD1	2.85	0.44
1:B:48:ARG:NH1	1:B:48:ARG:HB3	2.32	0.44
1:A:7:GLU:HG3	1:A:31:LEU:HD23	1.99	0.44
1:A:135:ASN:HB2	1:A:138:LEU:HD21	2.00	0.44
1:A:99:PRO:HG2	1:A:100:TYR:CE2	2.53	0.44
1:C:58:VAL:HG13	1:C:132:ILE:HG22	2.00	0.44
1:B:99:PRO:O	1:B:101:PRO:HD3	2.18	0.43
1:A:72:ASN:N	1:A:73:PRO:HD3	2.33	0.43
1:B:147:TRP:HB2	1:B:152:TYR:CE1	2.53	0.43
1:D:7:GLU:HG3	1:D:31:LEU:HD23	2.00	0.43
1:D:169:GLU:O	1:D:170:ASN:HB2	2.17	0.43
1:C:155:CYS:O	1:C:156:SER:HB2	2.18	0.43
1:A:11:LEU:HB3	1:A:114:PHE:HB2	2.00	0.43
1:D:91:GLN:O	1:D:94:ILE:HG22	2.19	0.43
3:D:504:4PT:O0H	3:D:504:4PT:OP1	2.36	0.43
1:A:35:HIS:CE1	1:D:35:HIS:HA	2.53	0.43
1:B:143:HIS:CD2	1:B:152:TYR:CE2	3.07	0.43
1:C:30:PHE:C	1:C:31:LEU:HD23	2.43	0.43
1:B:31:LEU:HD11	1:B:49:ARG:HH11	1.84	0.43
1:C:26:LYS:HG3	1:C:28:ARG:HH21	1.84	0.43
1:B:74:PRO:C	1:B:78:GLN:HG3	2.44	0.42
1:D:39:TYR:CZ	1:D:54:GLY:HA3	2.54	0.42
1:B:21:SER:HA	1:B:22:PRO:HD3	1.85	0.42
1:A:35:HIS:HE1	1:D:35:HIS:HA	1.85	0.42
1:C:46:ARG:HD2	1:C:48:ARG:HH21	1.85	0.42
1:A:92:ILE:CD1	1:B:9:ILE:HG13	2.49	0.42
1:A:142:TYR:CZ	1:A:144:PRO:HG3	2.55	0.42
1:C:11:LEU:HB2	1:C:98:PHE:CG	2.54	0.42
1:C:116:PRO:HG3	1:D:91:GLN:NE2	2.34	0.42
1:D:106:TYR:CE2	1:D:108:GLU:HB2	2.54	0.42
1:B:14:SER:HG	3:B:502:4PT:H6	1.83	0.42
1:B:40:TYR:CD1	1:B:40:TYR:N	2.86	0.42
1:B:155:CYS:HB3	1:B:165:CYS:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:HG23	1:B:102:PHE:HA	2.02	0.42
1:C:13:ARG:HG2	1:C:14:SER:H	1.82	0.42
1:C:17:LYS:HG2	1:C:108:GLU:HG2	2.01	0.42
1:A:21:SER:HA	1:A:22:PRO:HD3	1.73	0.42
1:A:68:VAL:HA	1:A:69:PRO:HD3	1.82	0.42
3:B:502:4PT:H0E2	3:B:502:4PT:C0P	2.45	0.41
1:C:23:LEU:N	1:C:23:LEU:HD23	2.35	0.41
1:A:52:LYS:CE	1:A:55:SER:HB2	2.50	0.41
1:B:143:HIS:HB3	1:B:154:CYS:SG	2.60	0.41
1:A:7:GLU:HA	1:A:30:PHE:O	2.19	0.41
1:B:152:TYR:HB3	1:B:154:CYS:SG	2.60	0.41
1:C:23:LEU:HG	1:C:23:LEU:O	2.19	0.41
1:C:110:PRO:HD3	1:C:160:LYS:HG2	2.01	0.41
1:C:138:LEU:O	1:C:140:GLN:NE2	2.54	0.41
1:C:148:ILE:HG22	1:C:149:ASP:N	2.34	0.41
1:D:79:ILE:H	1:D:79:ILE:HG13	1.56	0.41
1:A:35:HIS:HA	1:D:35:HIS:CE1	2.55	0.41
1:D:57:ASP:HB2	1:D:60:LYS:HG3	2.03	0.41
1:A:68:VAL:O	1:A:146:PHE:CE2	2.74	0.41
1:B:27:LYS:HE2	1:B:98:PHE:HE1	1.86	0.41
1:B:94:ILE:HG23	1:B:95:ILE:HG23	2.02	0.41
1:C:35:HIS:C	1:C:36:LYS:HG3	2.46	0.41
1:A:40:TYR:CD2	1:A:52:LYS:HA	2.56	0.41
1:B:130:ASN:HA	1:B:133:ARG:NE	2.36	0.41
1:C:7:GLU:HA	1:C:30:PHE:O	2.20	0.41
1:D:23:LEU:HD22	1:D:23:LEU:HA	1.90	0.41
1:D:147:TRP:CZ3	1:D:160:LYS:NZ	2.87	0.41
1:A:63:CYS:HB2	1:A:139:VAL:HG22	2.02	0.41
1:D:36:LYS:CA	1:D:58:VAL:HG23	2.48	0.41
1:D:64:VAL:HG21	1:D:128:LEU:HB3	2.03	0.41
1:A:64:VAL:HG21	1:A:128:LEU:HB3	2.03	0.40
1:B:147:TRP:CD2	1:B:160:LYS:HE3	2.56	0.40
1:C:24:ASN:OD1	1:C:24:ASN:C	2.64	0.40
1:A:67:VAL:O	1:A:68:VAL:C	2.64	0.40
1:B:148:ILE:HG22	1:B:149:ASP:CG	2.46	0.40
1:D:155:CYS:O	1:D:156:SER:HB2	2.22	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	154/169 (91%)	139 (90%)	15 (10%)	0	100	100
1	B	157/169 (93%)	141 (90%)	14 (9%)	2 (1%)	9	19
1	C	149/169 (88%)	130 (87%)	13 (9%)	6 (4%)	2	3
1	D	157/169 (93%)	147 (94%)	10 (6%)	0	100	100
All	All	617/676 (91%)	557 (90%)	52 (8%)	8 (1%)	9	19

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	19	LYS
1	C	16	GLN
1	C	107	ASP
1	C	108	GLU
1	C	150	GLY
1	B	107	ASP
1	C	148	ILE
1	C	99	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	148/158 (94%)	140 (95%)	8 (5%)	20	40
1	B	151/158 (96%)	143 (95%)	8 (5%)	20	41

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	145/158 (92%)	141 (97%)	4 (3%)	38	64
1	D	151/158 (96%)	146 (97%)	5 (3%)	33	59
All	All	595/632 (94%)	570 (96%)	25 (4%)	26	50

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LYS
1	A	40	TYR
1	A	67	VAL
1	A	70	GLU
1	A	97	ARG
1	A	140	GLN
1	A	151	GLN
1	A	168	LEU
1	B	14	SER
1	B	20	THR
1	B	40	TYR
1	B	71	LYS
1	B	101	PRO
1	B	107	ASP
1	B	137	ASP
1	B	166	GLN
1	C	40	TYR
1	C	108	GLU
1	C	137	ASP
1	C	158	THR
1	D	40	TYR
1	D	71	LYS
1	D	137	ASP
1	D	140	GLN
1	D	166	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	126	HIS
1	B	78	GLN
1	C	78	GLN

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Mol	Chain	Res	Type
1	C	126	HIS
1	C	130	ASN
1	C	140	GLN
1	D	78	GLN
1	D	127	GLN
1	D	130	ASN
1	D	151	GLN
1	D	170	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	4PT	C	503	-	28,28,43	1.58	3 (10%)	46,46,64	0.85	0
3	4PT	B	502	-	43,43,43	1.59	5 (11%)	58,64,64	1.25	5 (8%)
3	4PT	D	504	-	43,43,43	1.58	5 (11%)	58,64,64	1.12	3 (5%)
3	4PT	A	501	-	28,28,43	1.58	3 (10%)	46,46,64	0.81	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	4PT	C	503	-	-	1/20/44/65	0/1/1/1
3	4PT	B	502	-	2/2/13/13	11/41/65/65	0/1/1/1
3	4PT	D	504	-	2/2/13/13	9/41/65/65	0/1/1/1
3	4PT	A	501	-	-	4/20/44/65	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	4PT	C0I-C0G	-5.39	1.39	1.50
3	D	504	4PT	C0I-C0G	-5.38	1.39	1.50
3	A	501	4PT	P1-OP2	3.56	1.61	1.50
3	C	503	4PT	P1-OP2	3.55	1.61	1.50
3	C	503	4PT	P3-O12	3.55	1.61	1.50
3	D	504	4PT	P4-OP4	3.54	1.61	1.50
3	A	501	4PT	P3-O12	3.54	1.61	1.50
3	B	502	4PT	P4-OP4	3.54	1.61	1.50
3	C	503	4PT	P5-OP9	3.53	1.61	1.50
3	A	501	4PT	P5-OP9	3.52	1.61	1.50
3	B	502	4PT	P5-OP9	3.52	1.61	1.50
3	B	502	4PT	P3-O12	3.51	1.61	1.50
3	D	504	4PT	P5-OP9	3.50	1.61	1.50
3	D	504	4PT	P3-O12	3.49	1.61	1.50
3	B	502	4PT	C0J-C0I	-2.25	1.39	1.51
3	D	504	4PT	C0J-C0I	-2.24	1.39	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	504	4PT	O0H-C0G-C0I	3.65	119.33	109.37
3	B	502	4PT	O0H-C0G-C0I	3.47	118.85	109.37
3	D	504	4PT	C0J-C0I-C0G	3.41	119.73	113.76
3	B	502	4PT	C0J-C0I-C0G	3.08	119.17	113.76
3	D	504	4PT	O0F-C0L-C0N	3.00	120.97	111.83
3	B	502	4PT	C0B-C0C-C0E	-2.83	105.19	111.78
3	B	502	4PT	O0F-C0L-C0N	2.82	120.45	111.83
3	B	502	4PT	O0D-C0G-C0I	2.48	121.55	110.27

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	502	4PT	C0G
3	B	502	4PT	C0C
3	D	504	4PT	C0G
3	D	504	4PT	C0C

All (25) torsion outliers are listed below:

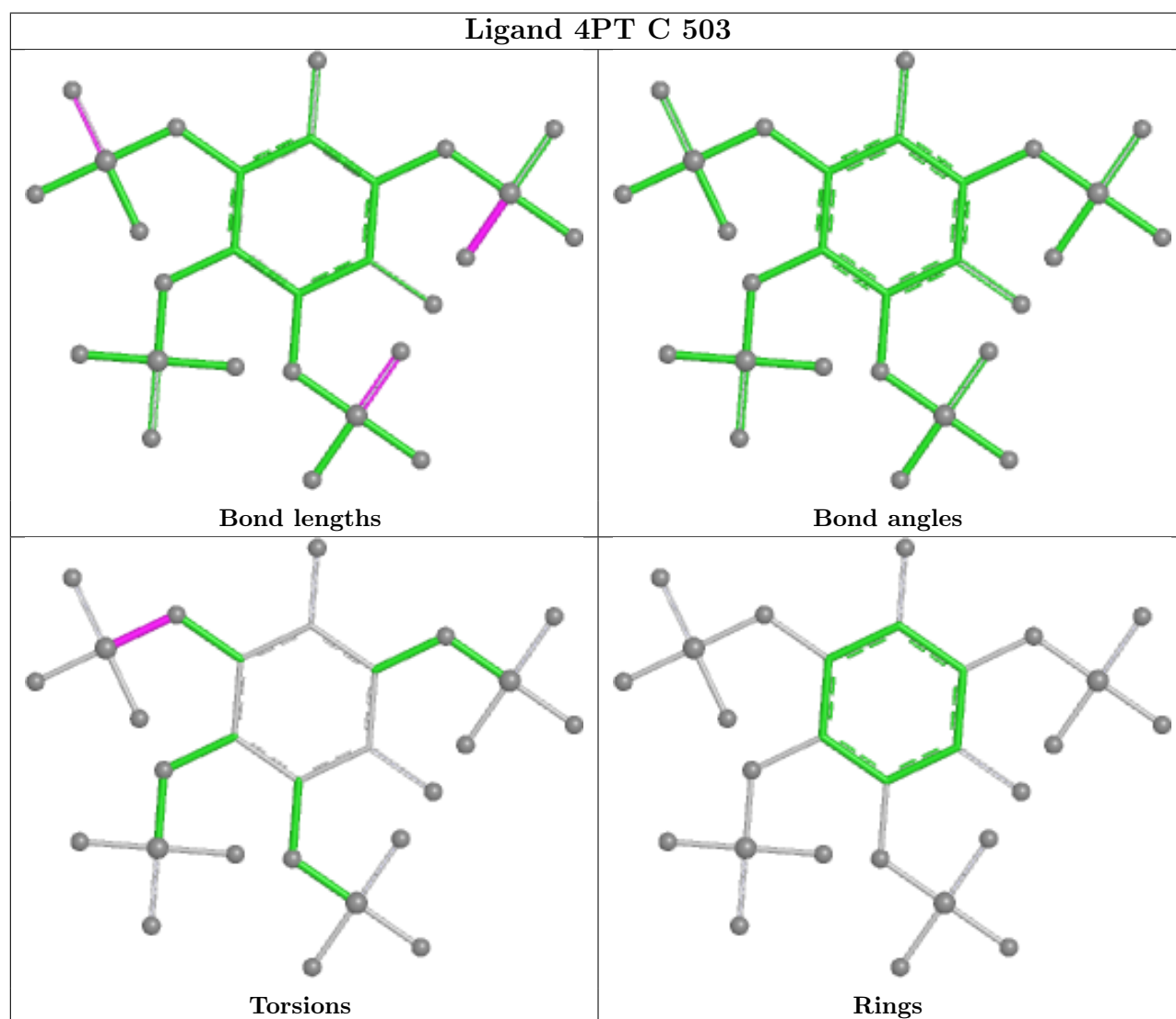
Mol	Chain	Res	Type	Atoms
3	B	502	4PT	O0M-C0L-O0F-C0E
3	B	502	4PT	C0N-C0L-O0F-C0E
3	B	502	4PT	O0H-C0G-O0D-C0C
3	B	502	4PT	C0I-C0G-O0D-C0C
3	B	502	4PT	O0H-C0G-C0I-C0J
3	D	504	4PT	O0M-C0L-O0F-C0E
3	D	504	4PT	C0N-C0L-O0F-C0E
3	D	504	4PT	O0H-C0G-O0D-C0C
3	D	504	4PT	O0D-C0G-C0I-C0J
3	A	501	4PT	C4-C5-O5-P5
3	B	502	4PT	C1-O1-P1-OP3
3	A	501	4PT	C6-C5-O5-P5
3	B	502	4PT	C2-C1-O1-P1
3	D	504	4PT	C0E-C0C-O0D-C0G
3	B	502	4PT	C6-C1-O1-P1
3	D	504	4PT	O0H-C0G-C0I-C0J
3	D	504	4PT	O0F-C0L-C0N-C0P
3	B	502	4PT	C1-O1-P1-OP2
3	A	501	4PT	C4-O4-P4-OP5
3	A	501	4PT	C4-O4-P4-OP6
3	C	503	4PT	C3-O3-P3-O10
3	B	502	4PT	O0F-C0L-C0N-C0P
3	D	504	4PT	C4-O4-P4-OP4
3	D	504	4PT	O0M-C0L-C0N-C0P
3	B	502	4PT	O0M-C0L-C0N-C0P

There are no ring outliers.

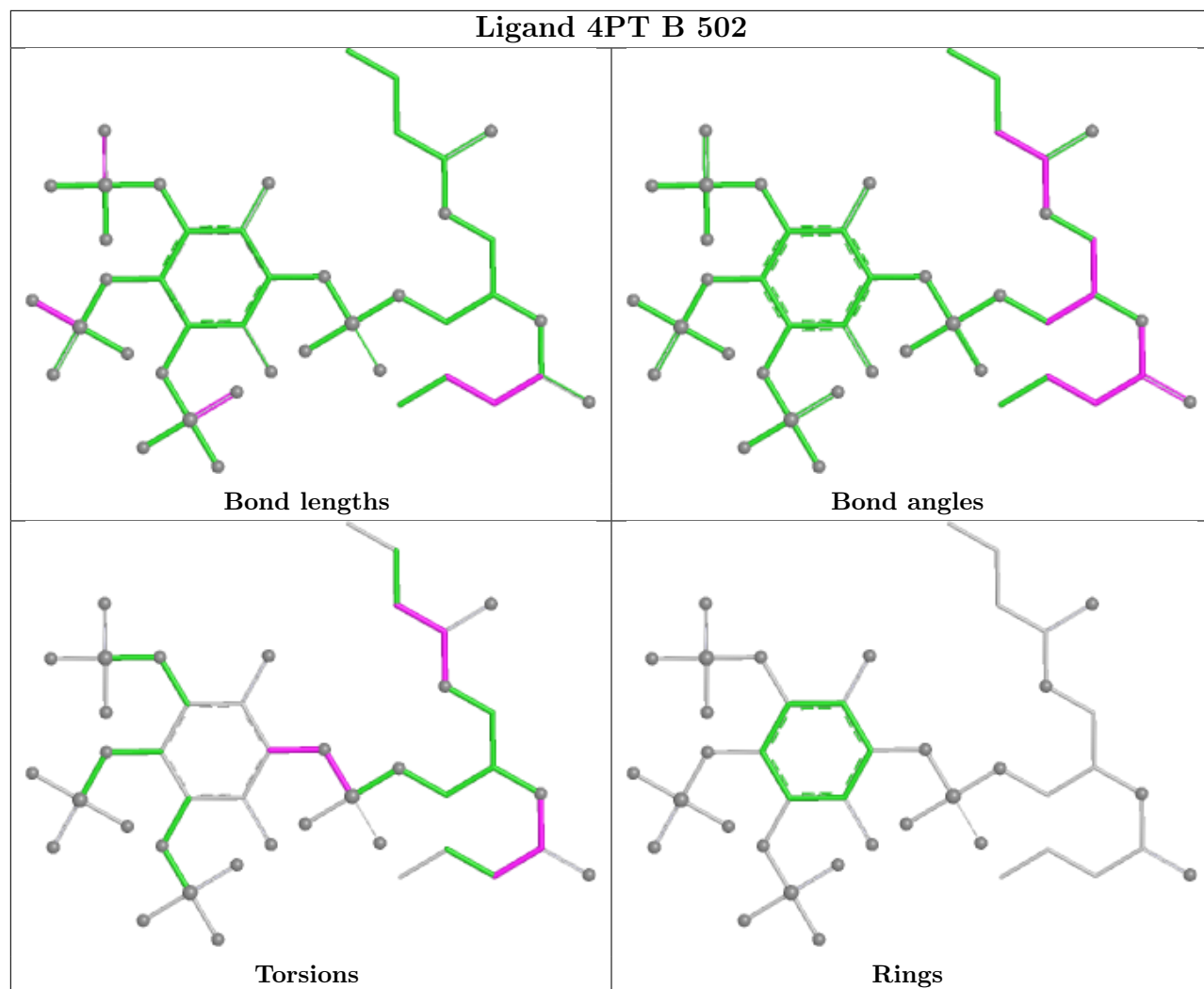
4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	4PT	1	0
3	B	502	4PT	5	0
3	D	504	4PT	4	0
3	A	501	4PT	1	0

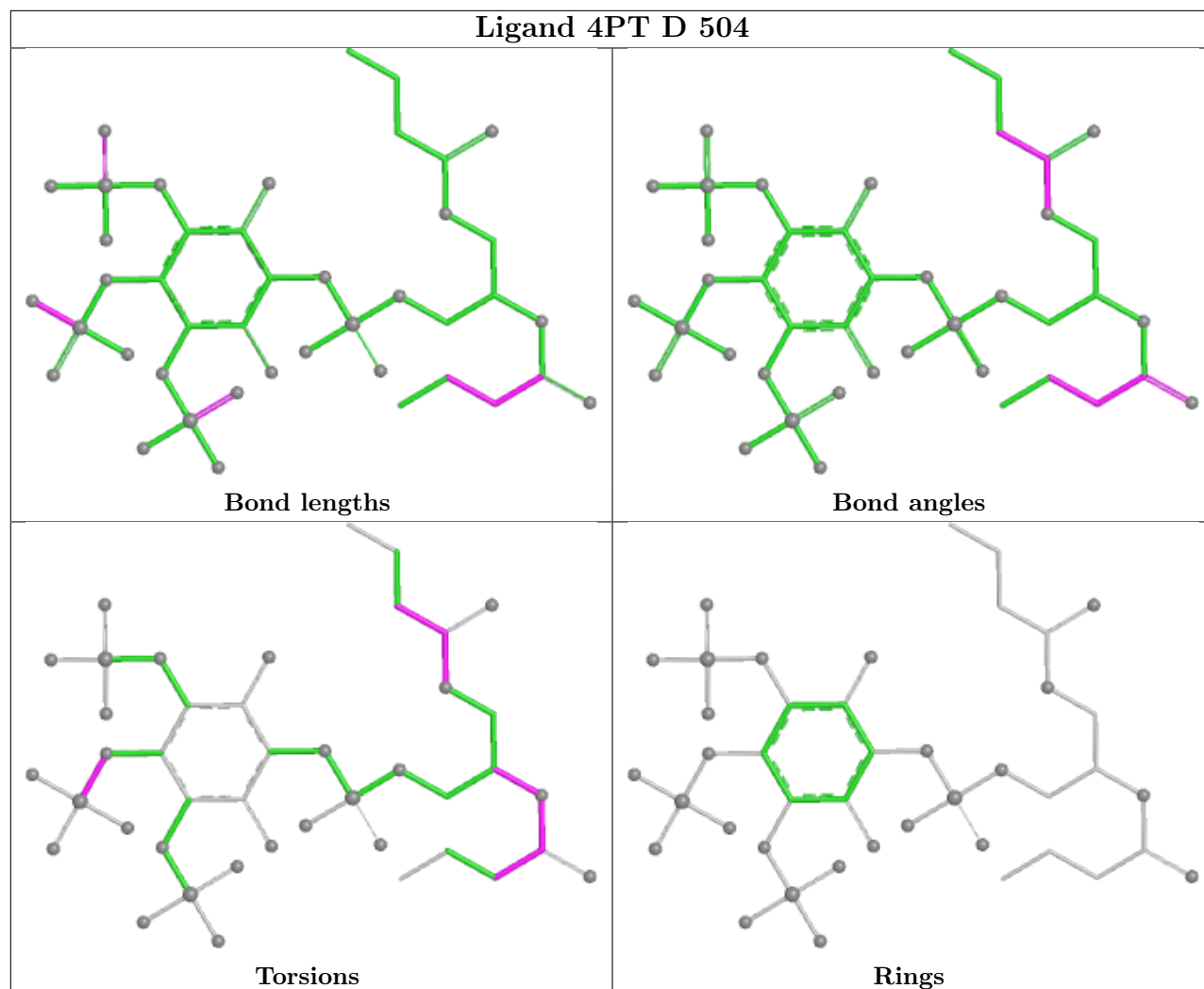
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

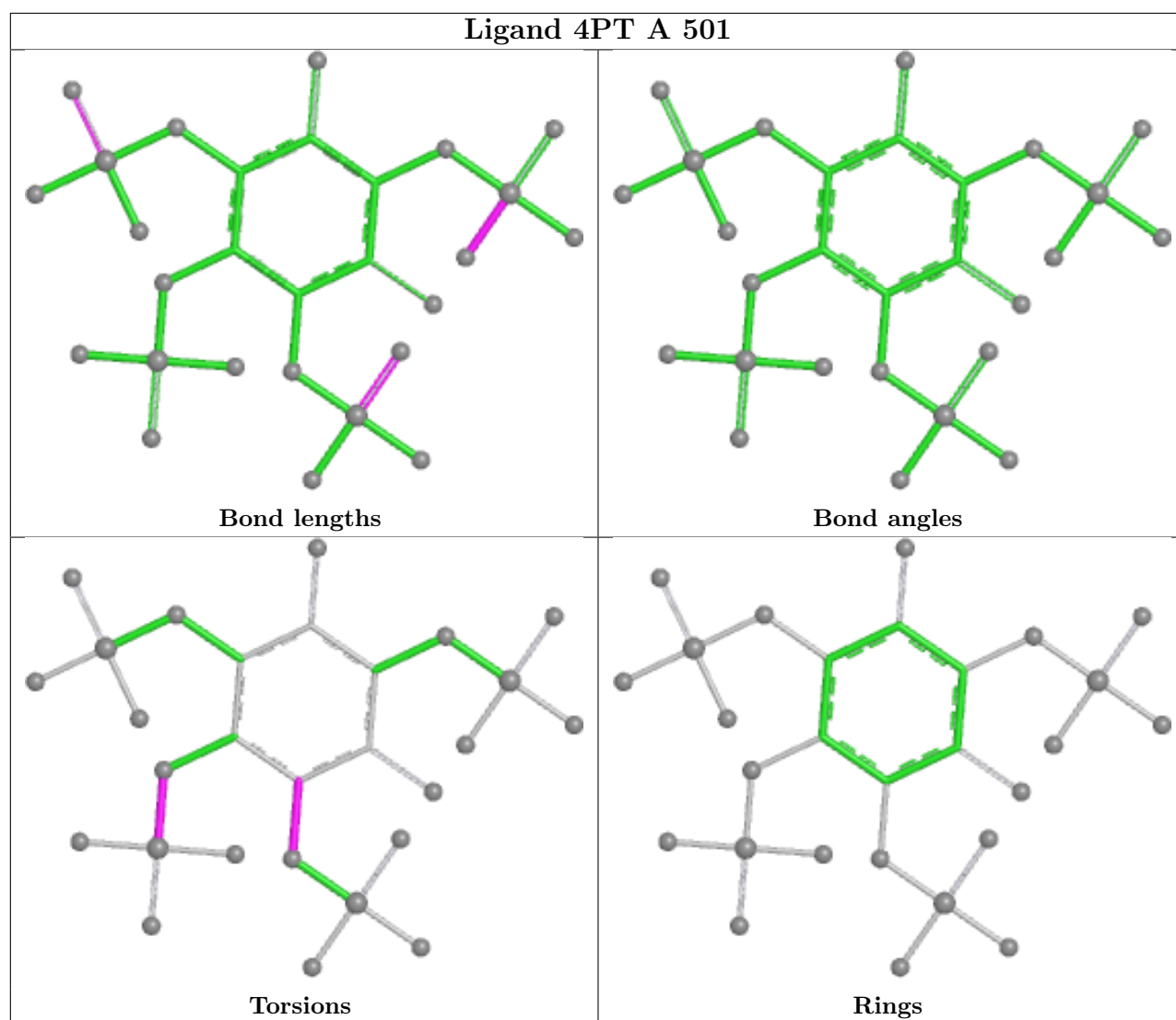


Ligand 4PT B 502



Ligand 4PT D 504





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	158/169 (93%)	0.63	10 (6%) 26 21	31, 50, 81, 88	0
1	B	161/169 (95%)	0.47	14 (8%) 16 13	24, 47, 77, 91	0
1	C	155/169 (91%)	0.86	12 (7%) 19 16	36, 59, 85, 99	0
1	D	161/169 (95%)	0.19	5 (3%) 51 46	23, 44, 65, 92	0
All	All	635/676 (93%)	0.53	41 (6%) 25 20	23, 50, 80, 99	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	14	SER	3.9
1	A	150	GLY	3.6
1	B	21	SER	3.5
1	B	152	TYR	3.4
1	A	153	LEU	3.0
1	C	109	GLY	3.0
1	B	147	TRP	2.8
1	A	73	PRO	2.7
1	C	148	ILE	2.7
1	C	164	GLY	2.6
1	A	155	CYS	2.6
1	B	140	GLN	2.6
1	C	134	TYR	2.6
1	A	75	PRO	2.6
1	B	109	GLY	2.5
1	B	20	THR	2.5
1	C	50	GLY	2.5
1	B	155	CYS	2.5
1	C	154	CYS	2.5
1	C	54	GLY	2.4
1	C	147	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	167	ILE	2.4
1	A	20	THR	2.3
1	D	79	ILE	2.3
1	C	150	GLY	2.3
1	B	19	LYS	2.2
1	C	70	GLU	2.2
1	A	89	MET	2.2
1	D	148	ILE	2.2
1	B	137	ASP	2.1
1	A	147	TRP	2.1
1	B	159	ALA	2.1
1	A	72	ASN	2.1
1	B	133	ARG	2.1
1	B	170	ASN	2.1
1	B	90	GLU	2.1
1	A	149	ASP	2.1
1	C	89	MET	2.1
1	D	168	LEU	2.0
1	D	80	PRO	2.0
1	B	89	MET	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	4PT	C	503	28/43	0.85	0.13	103,108,110,111	0
3	4PT	B	502	43/43	0.90	0.16	42,62,76,77	0
3	4PT	A	501	28/43	0.91	0.11	60,71,77,79	0

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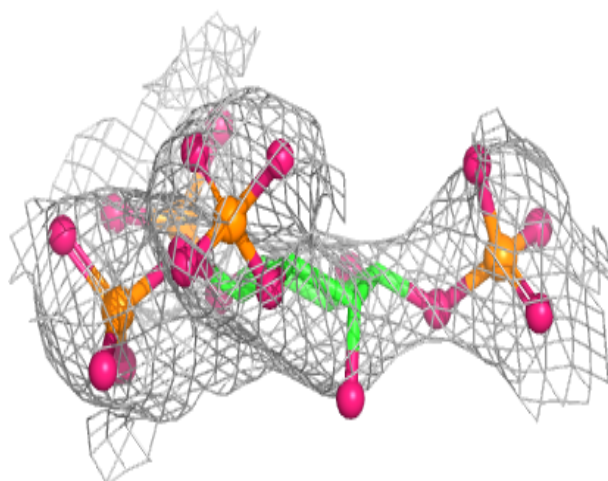
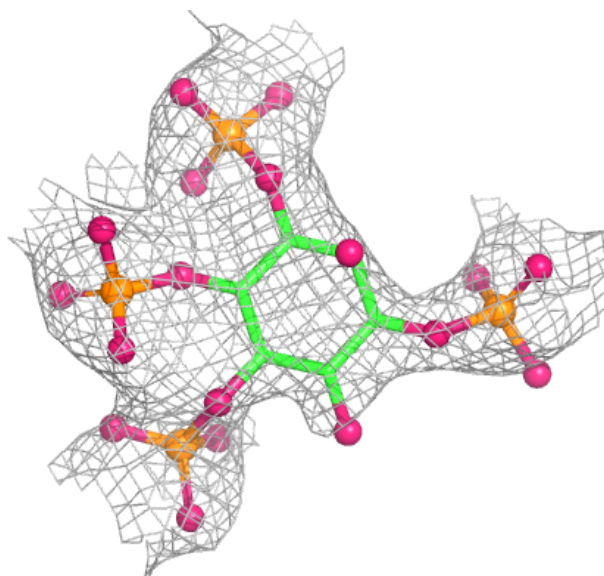
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	4PT	D	504	43/43	0.93	0.12	33,47,84,86	0
2	ZN	A	301	1/1	0.98	0.04	44,44,44,44	0
2	ZN	C	303	1/1	0.98	0.04	55,55,55,55	0
2	ZN	B	302	1/1	0.99	0.04	42,42,42,42	0
2	ZN	D	304	1/1	1.00	0.03	36,36,36,36	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.

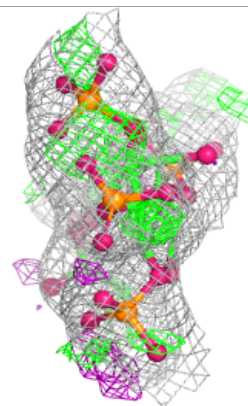
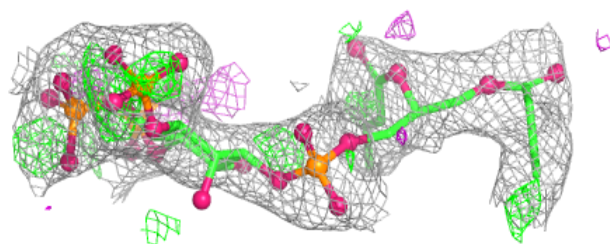
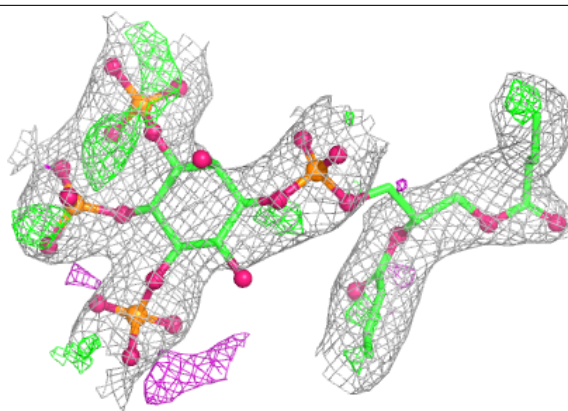
Electron density around 4PT C 503:

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



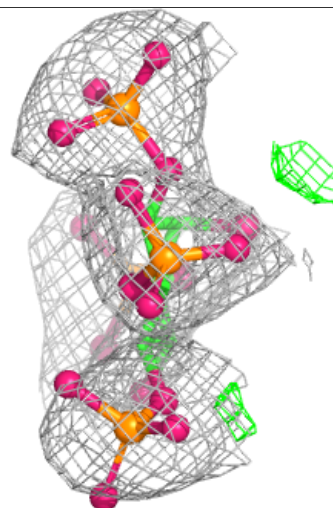
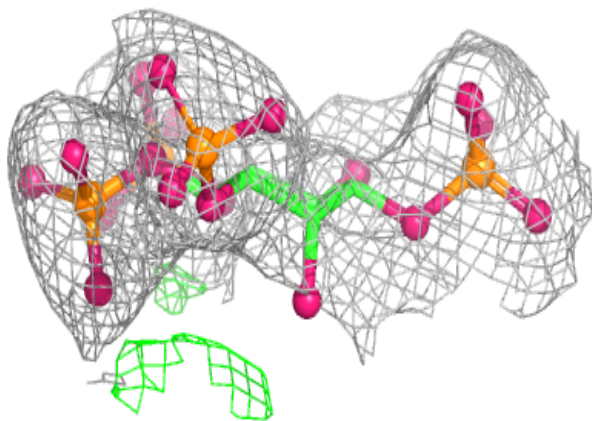
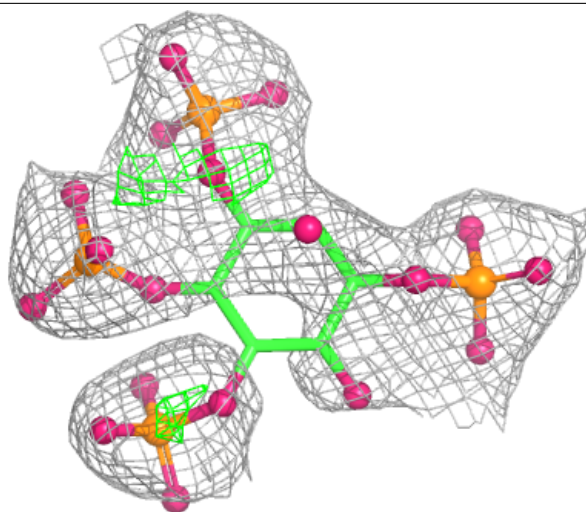
Electron density around 4PT B 502:

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



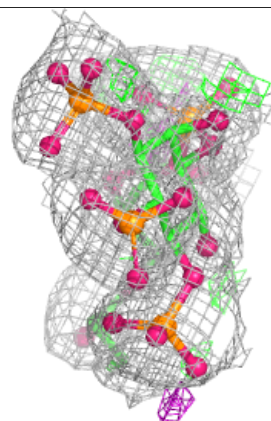
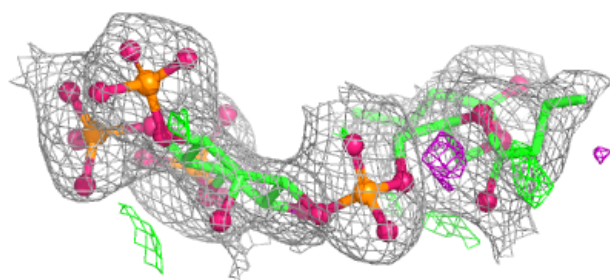
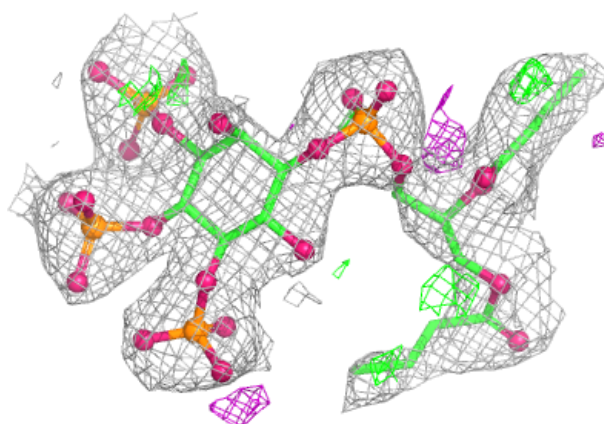
Electron density around 4PT A 501:

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



Electron density around 4PT D 504:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.