



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 12:06 AM UTC

PDB ID : 6YVY / pdb_00006yvy
Title : FOCAL ADHESION KINASE CATALYTIC DOMAIN IN COMPLEX WITH
4-{[4-{[(1R,2R)-2-(dimethylamino)cyclopentyl]amino}-5-(trifluoromethyl)pyr
imidin-2-yl]amino}-N-methylbenzenesulfonamide
Authors : Musil, D.; Amaral, M.
Deposited on : 2020-04-28
Resolution : 1.92 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

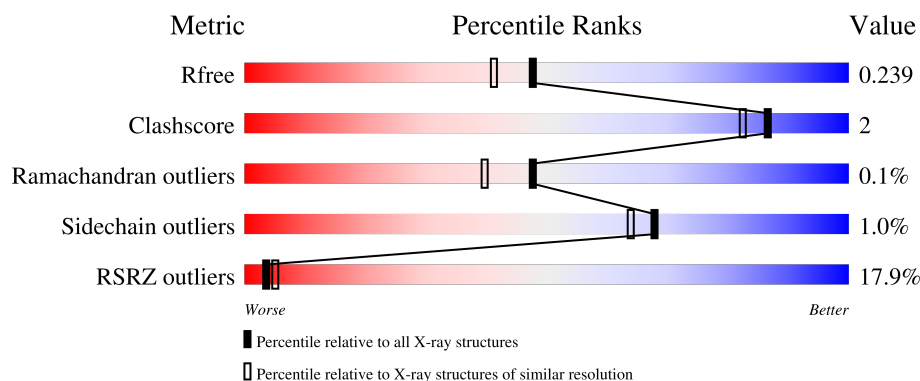
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	282	9% 87% 7% 6%
1	B	282	12% 84% 7% 9%
1	C	282	9% 88% • 9%
1	D	282	35% 86% • 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9023 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.

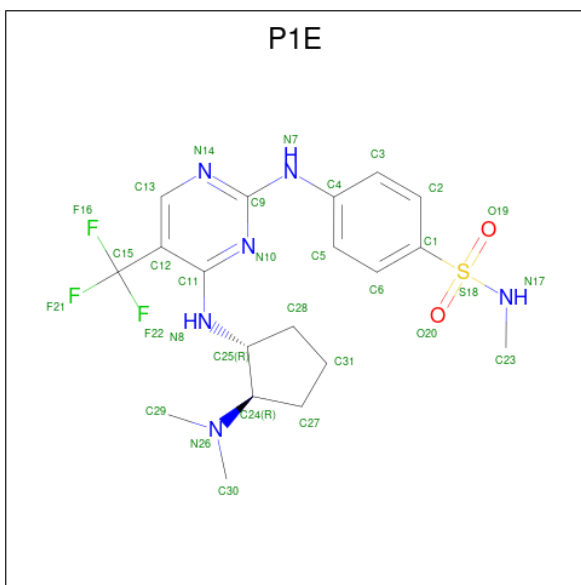
- Molecule 1 is a protein called Focal adhesion kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	3	0
			2154	1372	376	388	18			
1	B	257	Total	C	N	O	S	0	2	0
			2081	1330	361	373	17			
1	C	258	Total	C	N	O	S	0	3	0
			2099	1336	366	380	17			
1	D	255	Total	C	N	O	S	0	0	0
			2053	1313	355	368	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	408	GLY	-	expression tag	UNP Q05397
A	409	SER	-	expression tag	UNP Q05397
A	410	GLY	-	expression tag	UNP Q05397
B	408	GLY	-	expression tag	UNP Q05397
B	409	SER	-	expression tag	UNP Q05397
B	410	GLY	-	expression tag	UNP Q05397
C	408	GLY	-	expression tag	UNP Q05397
C	409	SER	-	expression tag	UNP Q05397
C	410	GLY	-	expression tag	UNP Q05397
D	408	GLY	-	expression tag	UNP Q05397
D	409	SER	-	expression tag	UNP Q05397
D	410	GLY	-	expression tag	UNP Q05397

- Molecule 2 is 4-{[4-{[(1R,2R)-2-(dimethylamino)cyclopentyl]amino}-5-(trifluoromethyl)pyrimidin-2-yl]amino}-N-methylbenzenesulfonamide (CCD ID: P1E) (formula: C₁₉H₂₅F₃N₆O₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			31	19	3	6	2	1		
2	B	1	Total	C	F	N	O	S	0	0
			31	19	3	6	2	1		
2	C	1	Total	C	F	N	O	S	0	0
			31	19	3	6	2	1		
2	D	1	Total	C	F	N	O	S	0	0
			31	19	3	6	2	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



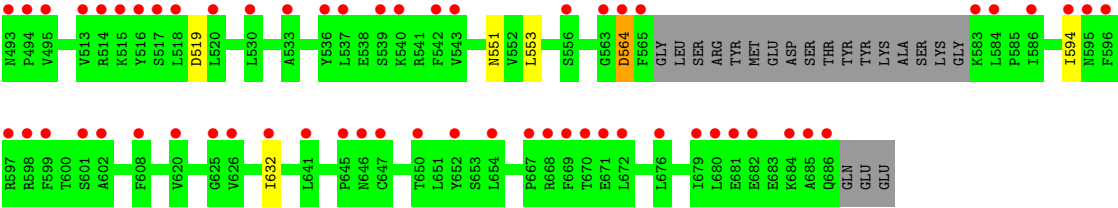
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	201	Total	O		0	0
			201	201			
4	B	133	Total	O		0	1
			134	134			
4	C	149	Total	O		0	0
			149	149			
4	D	23	Total	O		0	0
			23	23			

- Molecule 1: Focal adhesion kinase 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	113.97Å 75.86Å 172.64Å 90.00° 102.66° 90.00°	Depositor
Resolution (Å)	56.49 – 1.92 56.49 – 1.92	Depositor EDS
% Data completeness (in resolution range)	85.6 (56.49-1.92) 80.7 (56.49-1.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 1.92Å)	Xtriage
Refinement program	BUSTER 2.11.7 (19-MAR-2020)	Depositor
R, R_{free}	0.214 , 0.235 0.219 , 0.239	Depositor DCC
R_{free} test set	4419 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9023	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: SO4, P1E

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	0.81	1/2204 (0.0%)	1.02	2/2977 (0.1%)
1	B	0.73	0/2127	0.97	1/2876 (0.0%)
1	C	0.77	0/2145	1.00	2/2901 (0.1%)
1	D	0.60	0/2099	0.98	1/2838 (0.0%)
All	All	0.73	1/8575 (0.0%)	0.99	6/11592 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	643	MET	SD-CE	-8.13	1.59	1.79

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	519	ASP	CA-CB-CG	6.57	119.17	112.60
1	A	631	VAL	CA-C-N	6.26	128.68	120.60
1	A	631	VAL	C-N-CA	6.26	128.68	120.60
1	C	519	ASP	CA-CB-CG	5.46	118.06	112.60
1	C	438	GLN	N-CA-C	-5.17	102.83	110.48
1	D	519	ASP	CA-CB-CG	5.03	117.63	112.60

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	2154	0	2164	11	0
1	B	2081	0	2095	9	0
1	C	2099	0	2104	8	0
1	D	2053	0	2067	7	0
2	A	31	0	25	1	0
2	B	31	0	25	1	0
2	C	31	0	25	1	0
2	D	31	0	25	2	0
3	C	5	0	0	0	0
4	A	201	0	0	0	0
4	B	134	0	0	1	0
4	C	149	0	0	0	0
4	D	23	0	0	0	0
All	All	9023	0	8530	35	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 2.

All (35) 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:484:VAL:HG12	1:A:563:GLY:HA2	1.78	0.64
1:C:516:TYR:HB2	1:D:419:ARG:HD3	1.79	0.63
1:A:583:LYS:HE2	1:C:463:SER:HB2	1.81	0.61
1:C:516:TYR:HB2	1:D:419:ARG:CD	2.33	0.58
1:A:417:ILE:HD11	1:A:498:ILE:HD13	1.86	0.56
1:A:543:VAL:HG21	1:A:571:MET:HE2	1.87	0.56
1:A:484:VAL:CG1	1:A:563:GLY:HA2	2.36	0.55
1:C:671:GLU:HG3	1:C:675:GLN:HE21	1.72	0.54
1:D:594:ILE:HG21	1:D:632:ILE:HG12	1.91	0.52
1:B:484:VAL:HG21	1:B:553:LEU:HD12	1.92	0.52
1:D:471:GLU:O	1:D:475:MET:HG3	2.09	0.52
1:B:671:GLU:CG	1:B:675:GLN:HE21	2.22	0.51
1:B:598:ARG:NH1	4:B:802:HOH:O	2.37	0.51
1:A:415:TYR:HB2	1:A:476[B]:ARG:NH1	2.26	0.50
1:B:654:LEU:HD11	1:B:672:LEU:HD23	1.93	0.50
1:B:544:HIS:O	1:B:545[B]:ARG:HB2	2.12	0.49
1:A:632:ILE:H	1:A:632:ILE:HD12	1.78	0.48
2:D:701:P1E:H5	2:D:701:P1E:N10	2.29	0.48
1:B:514:ARG:CB	1:B:518:LEU:HG	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:654:LEU:HD11	1:A:672:LEU:HD23	1.97	0.46
1:D:553:LEU:HD21	2:D:701:P1E:H30B	1.99	0.44
1:C:654:LEU:HD11	1:C:672:LEU:HD23	1.99	0.43
2:C:802:P1E:H5	2:C:802:P1E:N10	2.34	0.43
1:B:484:VAL:HG13	1:B:500:GLU:HG2	2.01	0.42
1:D:551:ASN:ND2	1:D:564:ASP:OD2	2.52	0.42
1:A:457:LYS:HB2	1:A:457:LYS:HE3	1.62	0.42
1:B:555:SER:OG	1:B:561:LYS:HE3	2.20	0.42
1:B:417:ILE:HD11	1:B:487:ILE:HG22	2.00	0.42
2:B:701:P1E:H5	2:B:701:P1E:N10	2.35	0.41
2:A:701:P1E:H5	2:A:701:P1E:N10	2.36	0.41
1:C:671:GLU:CG	1:C:675:GLN:HE21	2.32	0.41
1:C:412:THR:O	1:C:412:THR:HG22	2.21	0.41
1:D:475:MET:HE1	1:D:484:VAL:HG23	2.03	0.41
1:A:484:VAL:HG23	1:A:500:GLU:HG2	2.02	0.40
1:A:597:ARG:HH11	1:C:411:SER:HB2	1.85	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	264/282 (94%)	260 (98%)	4 (2%)	0	100	100
1	B	255/282 (90%)	251 (98%)	4 (2%)	0	100	100
1	C	257/282 (91%)	253 (98%)	4 (2%)	0	100	100
1	D	251/282 (89%)	244 (97%)	6 (2%)	1 (0%)	30	19
All	All	1027/1128 (91%)	1008 (98%)	18 (2%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	564	ASP

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	238/249 (96%)	235 (99%)	3 (1%)	61	53
1	B	230/249 (92%)	228 (99%)	2 (1%)	70	67
1	C	233/249 (94%)	231 (99%)	2 (1%)	70	67
1	D	227/249 (91%)	225 (99%)	2 (1%)	70	67
All	All	928/996 (93%)	919 (99%)	9 (1%)	68	64

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	443	SER
1	A	461	SER
1	A	629	ASN
1	B	443	SER
1	B	629	ASN
1	C	412	THR
1	C	442	MET
1	D	443	SER
1	D	459	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	446	ASN
1	A	629	ASN
1	B	493	ASN
1	B	629	ASN
1	B	675	GLN
1	C	438	GLN
1	C	458	ASN

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Mol	Chain	Res	Type
1	C	675	GLN
1	D	438	GLN
1	D	470	GLN
1	D	493	ASN
1	D	551	ASN
1	D	618	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	P1E	D	701	-	32,33,33	0.23	0	46,49,49	0.65	1 (2%)
3	SO4	C	801	-	4,4,4	0.31	0	6,6,6	0.20	0
2	P1E	B	701	-	32,33,33	0.29	0	46,49,49	0.73	1 (2%)
2	P1E	C	802	-	32,33,33	0.37	0	46,49,49	0.74	1 (2%)
2	P1E	A	701	-	32,33,33	0.41	0	46,49,49	0.82	1 (2%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	P1E	A	701	-	-	1/27/37/37	0/3/3/3
2	P1E	B	701	-	-	1/27/37/37	0/3/3/3
2	P1E	D	701	-	-	1/27/37/37	0/3/3/3
2	P1E	C	802	-	-	1/27/37/37	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	802	P1E	C15-C12-C11	3.58	124.64	120.06
2	A	701	P1E	C15-C12-C11	3.51	124.55	120.06
2	B	701	P1E	C15-C12-C11	3.47	124.49	120.06
2	D	701	P1E	C15-C12-C11	3.44	124.45	120.06

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	P1E	C24-C25-N8-C11
2	B	701	P1E	C24-C25-N8-C11
2	C	802	P1E	C24-C25-N8-C11
2	D	701	P1E	C24-C25-N8-C11

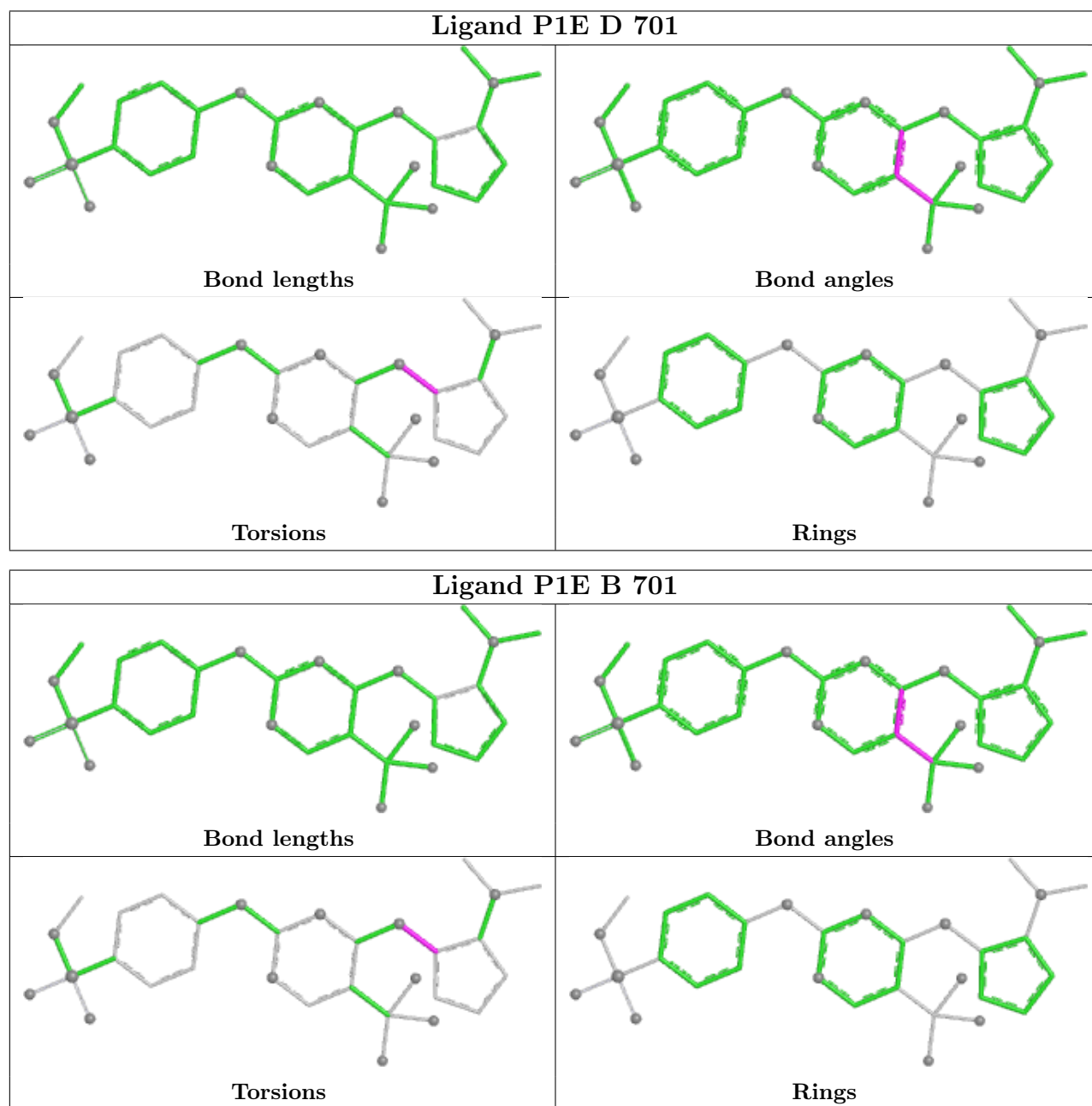
There are no ring outliers.

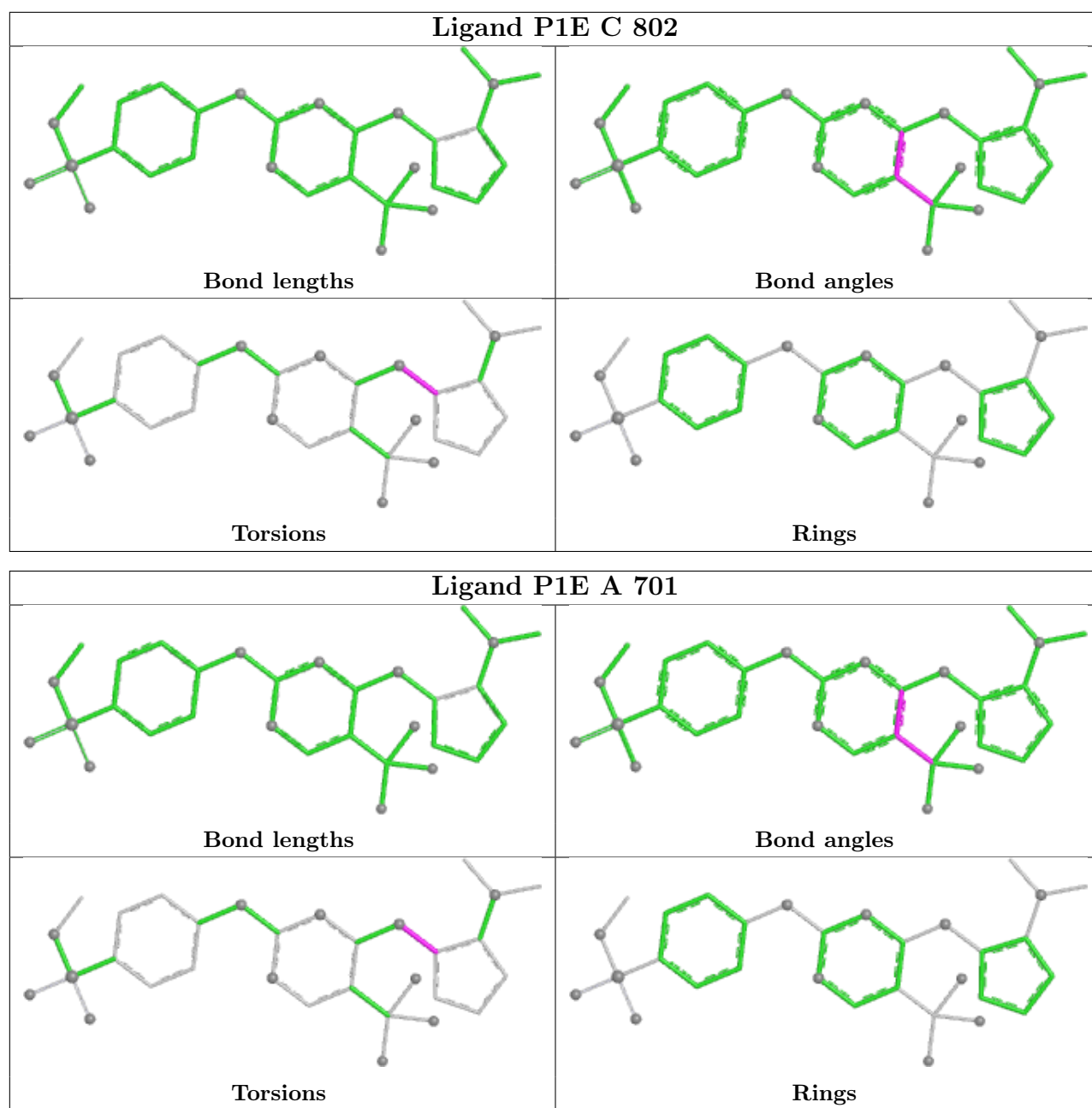
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	701	P1E	2	0
2	B	701	P1E	1	0
2	C	802	P1E	1	0
2	A	701	P1E	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.





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	265/282 (93%)	0.63	25 (9%) 14 17	12, 26, 46, 63	3 (1%)
1	B	257/282 (91%)	0.87	35 (13%) 7 9	13, 32, 62, 76	2 (0%)
1	C	258/282 (91%)	0.72	26 (10%) 12 16	12, 29, 53, 68	3 (1%)
1	D	255/282 (90%)	1.81	99 (38%) 1 0	31, 57, 94, 117	0
All	All	1035/1128 (91%)	1.00	185 (17%) 3 5	12, 34, 74, 117	8 (0%)

All (185) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	460	THR	5.6
1	A	516	TYR	5.3
1	D	494	PRO	4.9
1	B	459	CYS	4.9
1	D	459	CYS	4.9
1	D	493	ASN	4.9
1	C	444	PRO	4.9
1	D	415	TYR	4.6
1	D	516	TYR	4.6
1	B	567	LEU	4.5
1	C	459	CYS	4.5
1	D	680	LEU	4.4
1	D	491	THR	4.4
1	B	415	TYR	4.3
1	D	539	SER	4.3
1	D	469	LEU	4.2
1	D	464	VAL	4.2
1	D	517	SER	4.1
1	B	568	SER	4.1
1	D	565	PHE	4.1
1	D	465	ARG	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	571	MET	4.0
1	A	628	ASN	3.9
1	C	685	ALA	3.8
1	B	464	VAL	3.8
1	A	631	VAL	3.8
1	A	563	GLY	3.7
1	A	446	ASN	3.7
1	B	444	PRO	3.6
1	D	584	LEU	3.5
1	A	630	ASP	3.5
1	C	412	THR	3.5
1	D	518	LEU	3.5
1	A	445	GLU	3.5
1	D	684	LYS	3.5
1	D	492	GLU	3.4
1	C	465	ARG	3.4
1	D	542	PHE	3.4
1	C	584	LEU	3.4
1	A	412	THR	3.4
1	B	563	GLY	3.4
1	B	584	LEU	3.4
1	D	473	LEU	3.4
1	D	599	PHE	3.4
1	C	460	THR	3.3
1	B	417	ILE	3.3
1	D	458	ASN	3.3
1	D	672	LEU	3.3
1	D	475	MET	3.3
1	D	537	LEU	3.3
1	A	582	GLY	3.3
1	D	563	GLY	3.3
1	C	464	VAL	3.2
1	D	490	ILE	3.2
1	B	565	PHE	3.2
1	D	472	ALA	3.1
1	D	425	GLY	3.1
1	D	463	SER	3.1
1	B	466	GLU	3.1
1	C	491	THR	3.1
1	D	474	THR	3.1
1	A	447	PRO	3.1
1	B	461	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	685	ALA	3.1
1	D	632	ILE	3.1
1	D	686	GLN	3.1
1	D	433	PHE	3.1
1	A	685	ALA	3.1
1	D	564	ASP	3.0
1	D	626	VAL	3.0
1	D	625	GLY	3.0
1	D	457	LYS	3.0
1	D	646	ASN	3.0
1	B	422	ILE	3.0
1	D	520	LEU	2.9
1	B	686	GLN	2.9
1	D	417	ILE	2.9
1	B	463	SER	2.9
1	C	457	LYS	2.9
1	D	670	THR	2.9
1	D	456	CYS	2.9
1	B	635	ILE	2.9
1	C	462	ASP	2.9
1	A	686	GLN	2.9
1	B	462	ASP	2.9
1	A	570	TYR	2.8
1	D	645	PRO	2.8
1	B	518	LEU	2.8
1	C	473	LEU	2.8
1	D	487	ILE	2.8
1	B	457	LYS	2.8
1	D	540	LYS	2.8
1	D	468	PHE	2.8
1	D	601	SER	2.7
1	D	432	GLN	2.7
1	D	597	ARG	2.7
1	A	519	ASP	2.7
1	C	461	SER	2.7
1	C	415	TYR	2.7
1	D	543	VAL	2.7
1	D	536	TYR	2.6
1	A	413	ARG	2.6
1	B	596	PHE	2.6
1	C	413	ARG	2.6
1	A	459	CYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	676	LEU	2.6
1	C	442	MET	2.6
1	C	686	GLN	2.6
1	D	513	VAL	2.6
1	D	669	PHE	2.6
1	D	467	LYS	2.6
1	B	460	THR	2.6
1	D	596	PHE	2.5
1	A	632	ILE	2.5
1	B	446	ASN	2.5
1	C	411	SER	2.5
1	D	495	VAL	2.5
1	A	414	ASP	2.5
1	B	426	ARG	2.5
1	D	436	VAL	2.5
1	D	667	PRO	2.5
1	C	565	PHE	2.4
1	D	470	GLN	2.4
1	D	679	ILE	2.4
1	D	514	ARG	2.4
1	D	530	LEU	2.4
1	D	533	ALA	2.4
1	B	597	ARG	2.4
1	C	447	PRO	2.4
1	C	445	GLU	2.4
1	A	557	ASN	2.4
1	D	466	GLU	2.4
1	D	682	GLU	2.4
1	B	564	ASP	2.3
1	D	608	PHE	2.3
1	D	620	VAL	2.3
1	D	654	LEU	2.3
1	D	586	ILE	2.3
1	D	681	GLU	2.3
1	D	647	CYS	2.3
1	D	478	PHE	2.3
1	D	477	GLN	2.3
1	A	424	LEU	2.3
1	D	424	LEU	2.3
1	D	485	LYS	2.3
1	D	583	LYS	2.3
1	A	569	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	685	ALA	2.2
1	D	431	GLY	2.2
1	D	602	ALA	2.2
1	D	598	ARG	2.2
1	D	556	SER	2.2
1	A	684	LYS	2.2
1	C	518	LEU	2.2
1	C	463	SER	2.2
1	D	444	PRO	2.2
1	D	652	TYR	2.2
1	A	583	LYS	2.2
1	D	515	LYS	2.2
1	B	445	GLU	2.2
1	D	430	GLU	2.2
1	B	513	VAL	2.2
1	D	641	LEU	2.2
1	D	476	ARG	2.2
1	B	470	GLN	2.1
1	B	430	GLU	2.1
1	D	594	ILE	2.1
1	D	461	SER	2.1
1	C	423	GLU	2.1
1	B	448	ALA	2.1
1	B	443	SER	2.0
1	B	594	ILE	2.0
1	C	684	LYS	2.0
1	D	471	GLU	2.0
1	D	671	GLU	2.0
1	D	595	ASN	2.0
1	A	626	VAL	2.0
1	B	631	VAL	2.0
1	B	519	ASP	2.0
1	D	668	ARG	2.0
1	D	445	GLU	2.0
1	C	443[A]	SER	2.0
1	D	650	THR	2.0
1	D	446	ASN	2.0
1	D	484	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

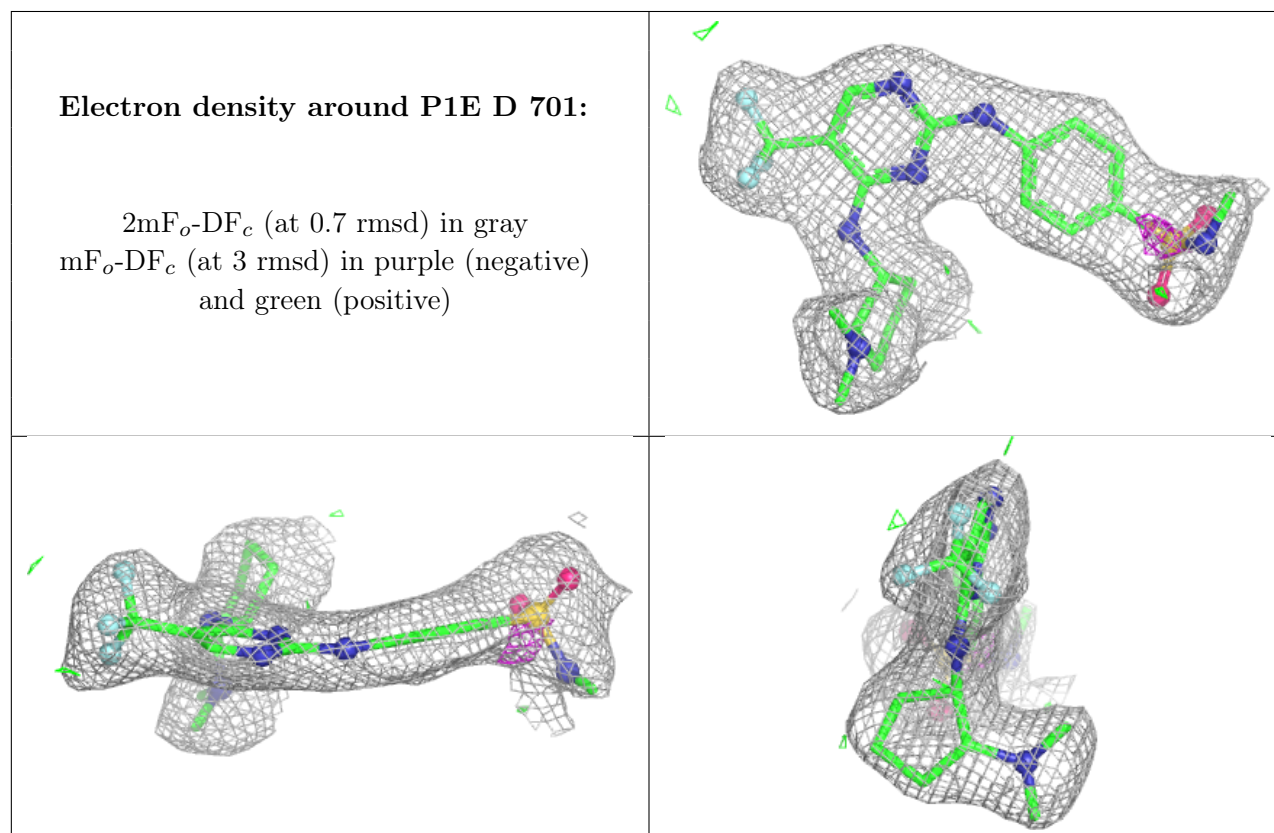
There are no oligosaccharides in this entry.

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.

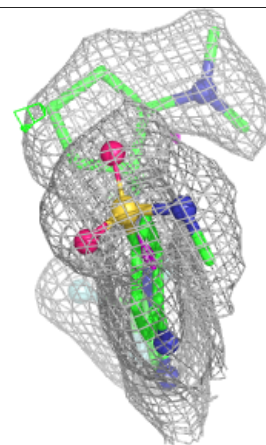
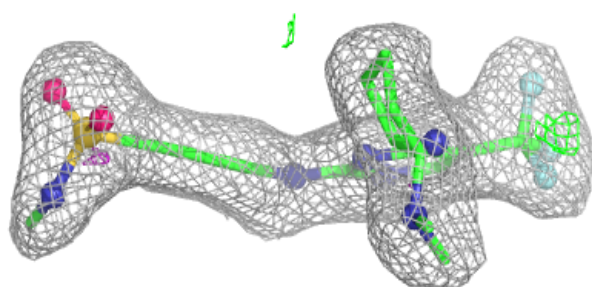
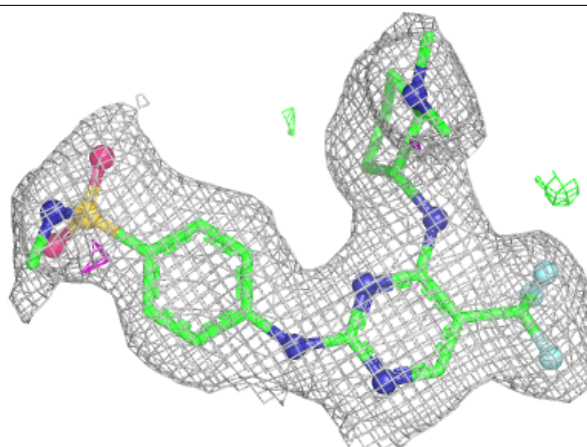
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	P1E	D	701	31/31	0.92	0.11	40,42,52,52	0
3	SO4	C	801	5/5	0.92	0.10	64,64,64,64	0
2	P1E	B	701	31/31	0.93	0.10	32,36,42,43	0
2	P1E	A	701	31/31	0.94	0.09	21,25,30,32	0
2	P1E	C	802	31/31	0.95	0.09	22,26,35,35	0

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.



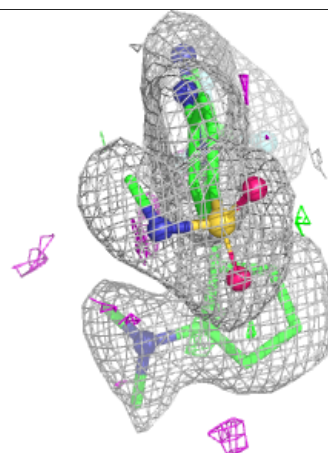
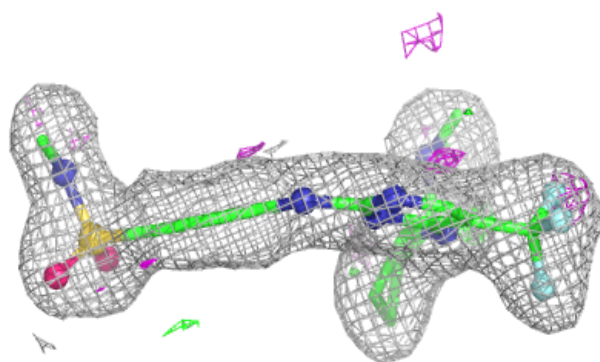
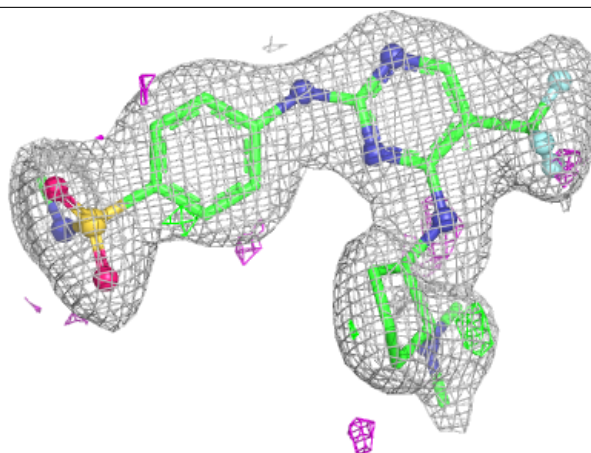
Electron density around P1E B 701:

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

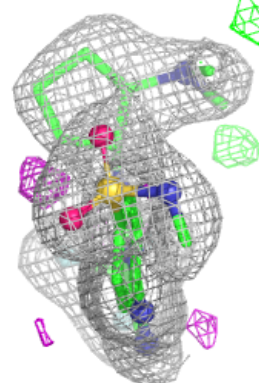
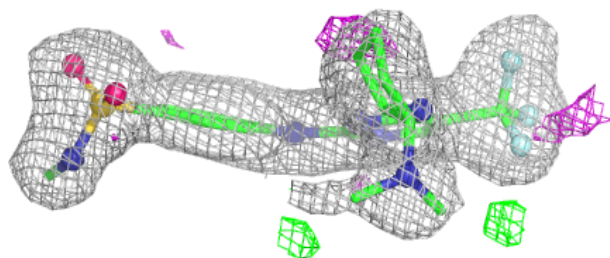
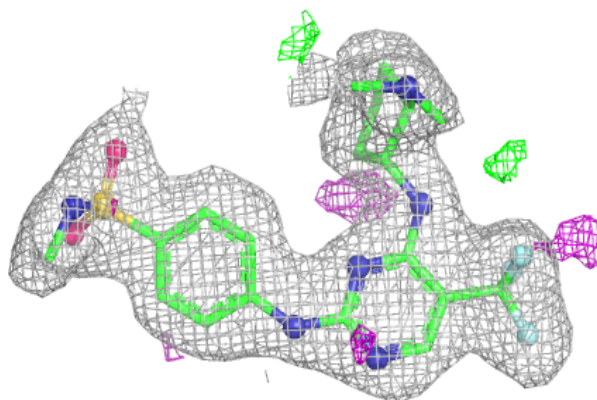


Electron density around P1E A 701:

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

**Electron density around P1E C 802:**

$2mF_o-DF_c$ (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.