



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:03 PM UTC

PDB ID : 3FX6 / pdb_00003fx6
Title : X-RAY crystallographic studies on the complex of carboxypeptidase A with the inhibitor using alpha-nitro ketone as the zinc-binding group
Authors : Wang, S.F.; Jin, J.Y.; Tian, G.R.
Deposited on : 2009-01-20
Resolution : 1.85 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

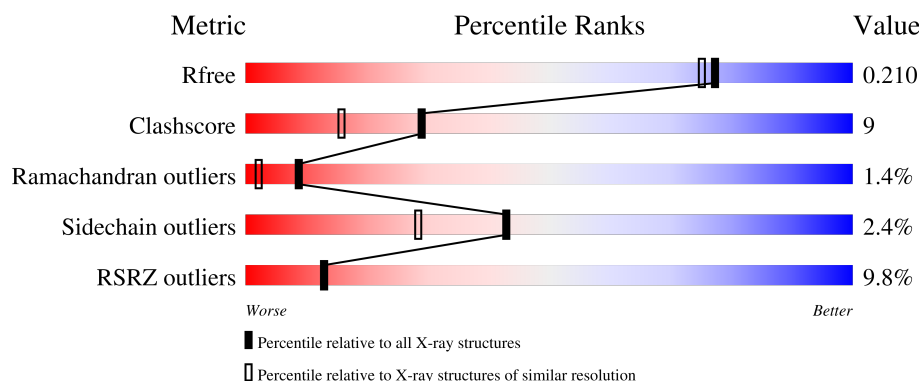
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	307	8% 79% 18% ..
1	C	307	12% 80% 16% ..
1	E	307	9% 80% 17% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7924 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 Carboxypeptidase A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	S	0	0	0
			2440	1563	403	469	5			
1	C	307	Total	C	N	O	S	0	0	0
			2440	1563	403	469	5			
1	E	307	Total	C	N	O	S	0	0	0
			2440	1563	403	469	5			

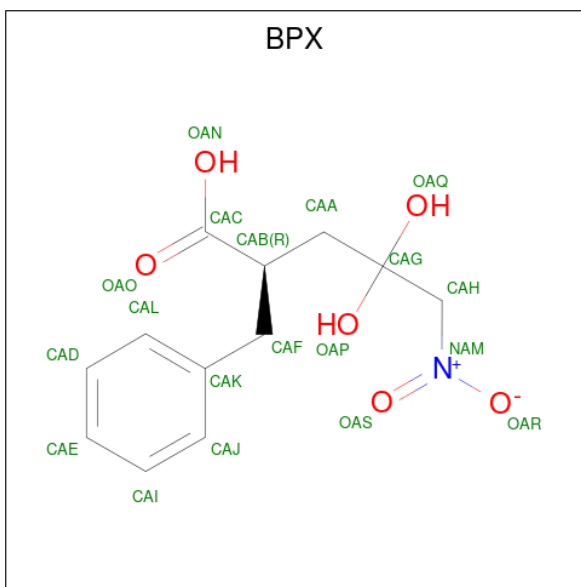
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	305	VAL	LEU	conflict	UNP P00730
C	305	VAL	LEU	conflict	UNP P00730
E	305	VAL	LEU	conflict	UNP P00730

- 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	C	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (2R)-4,4-dihydroxy-5-nitro-2-(phenylmethyl)pentanoic acid (CCD ID: BPX) (formula: C₁₂H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			19	12	1	6		
3	C	1	Total	C	N	O	0	0
			19	12	1	6		
3	E	1	Total	C	N	O	0	0
			19	12	1	6		

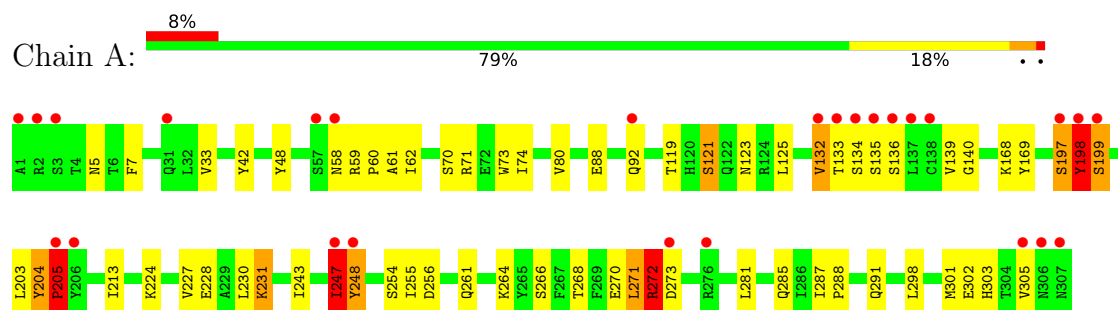
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	183	Total	O	0	0
			183	183		
4	C	176	Total	O	0	0
			176	176		
4	E	185	Total	O	0	0
			185	185		

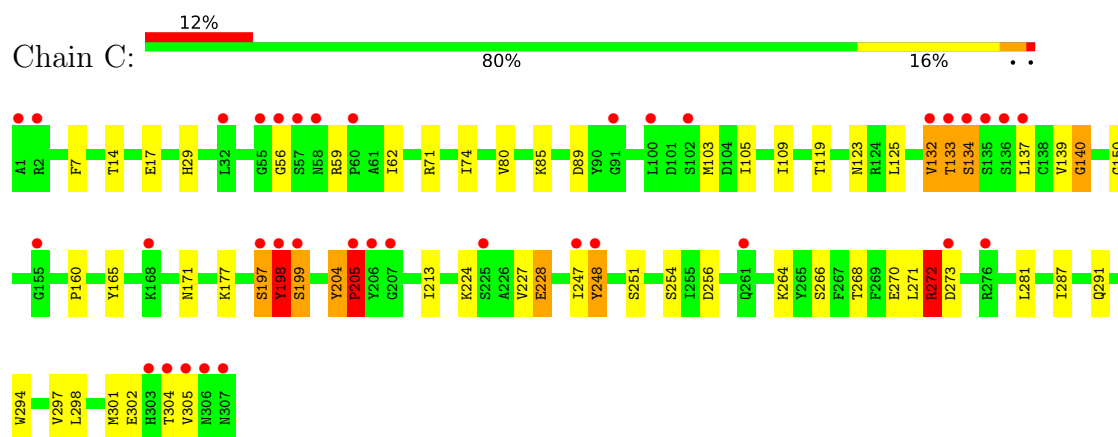
3 Residue-property plots [i](#)

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.

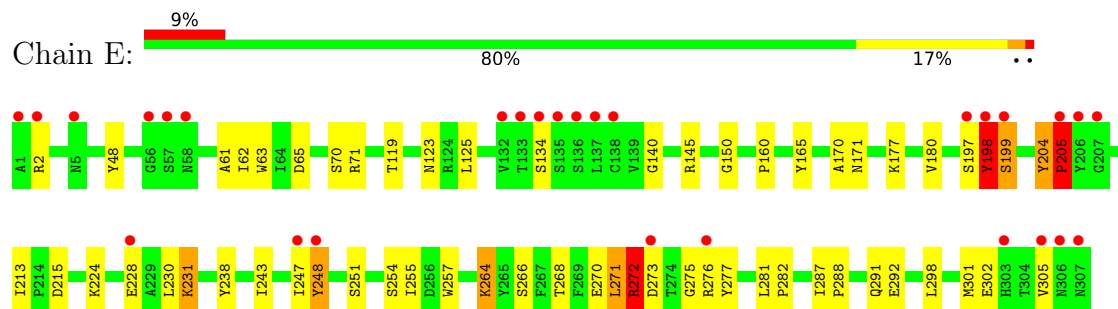
• Molecule 1: Carboxypeptidase A1



• Molecule 1: Carboxypeptidase A1



• Molecule 1: Carboxypeptidase A1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.36Å 59.75Å 99.51Å 90.00° 104.04° 90.00°	Depositor
Resolution (Å)	36.37 – 1.85 36.37 – 1.85	Depositor EDS
% Data completeness (in resolution range)	89.9 (36.37-1.85) 89.9 (36.37-1.85)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	22.33 (at 1.85Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.239 0.213 , 0.210	Depositor DCC
R_{free} test set	3217 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7924	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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 71.57 % 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 2.5664e-06. 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: BPX, 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	0.37	0/2506	0.98	19/3407 (0.6%)
1	C	0.37	0/2506	0.94	14/3407 (0.4%)
1	E	0.36	0/2506	0.94	12/3407 (0.4%)
All	All	0.36	0/7518	0.95	45/10221 (0.4%)

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	70	SER	N-CA-C	8.21	121.27	111.33
1	A	70	SER	N-CA-C	7.96	120.97	111.33
1	E	204	TYR	CA-C-N	7.79	129.57	119.84
1	E	204	TYR	C-N-CA	7.79	129.57	119.84
1	C	272	ARG	N-CA-C	7.78	127.37	110.80
1	E	197	SER	N-CA-C	7.56	121.87	109.24
1	A	197	SER	N-CA-C	7.50	121.76	109.24
1	E	271	LEU	N-CA-C	7.43	119.40	110.19
1	C	197	SER	N-CA-C	7.30	121.43	109.24
1	A	204	TYR	CA-C-N	7.13	128.76	119.84
1	A	204	TYR	C-N-CA	7.13	128.76	119.84
1	C	264	LYS	N-CA-C	7.10	118.67	111.07
1	E	140	GLY	N-CA-C	7.01	121.51	112.68
1	C	140	GLY	N-CA-C	6.81	121.50	112.65
1	A	272	ARG	N-CA-C	6.71	125.09	110.80
1	A	271	LEU	N-CA-C	6.60	119.70	109.60
1	E	272	ARG	N-CA-C	6.46	124.56	110.80
1	C	204	TYR	CA-C-N	6.36	127.79	119.84
1	C	204	TYR	C-N-CA	6.36	127.79	119.84
1	A	264	LYS	N-CA-C	6.28	118.64	111.11
1	A	205	PRO	N-CA-C	6.22	125.28	112.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	256	ASP	N-CA-C	-6.20	104.61	111.36
1	A	256	ASP	N-CA-C	-5.85	104.99	111.36
1	E	264	LYS	N-CA-C	5.79	118.05	111.11
1	E	205	PRO	N-CA-C	5.71	124.23	112.47
1	A	247	ILE	CB-CA-C	-5.66	106.90	111.71
1	A	254	SER	N-CA-C	5.65	117.44	111.28
1	A	61	ALA	N-CA-C	5.60	118.20	109.52
1	E	61	ALA	N-CA-C	5.52	118.07	109.52
1	C	198	TYR	N-CA-C	5.50	122.51	110.80
1	A	198	TYR	N-CA-C	5.50	122.51	110.80
1	C	205	PRO	N-CA-C	5.49	123.77	112.47
1	E	254	SER	N-CA-C	5.36	116.81	111.07
1	C	254	SER	N-CA-C	5.28	116.72	111.07
1	A	261	GLN	N-CA-C	-5.25	106.88	113.28
1	C	29	HIS	CA-C-N	5.24	125.05	119.28
1	C	29	HIS	C-N-CA	5.24	125.05	119.28
1	A	33	VAL	N-CA-C	5.21	115.98	108.48
1	A	121	SER	N-CA-C	5.17	119.48	112.04
1	A	73	TRP	N-CA-C	5.15	118.35	111.75
1	C	89	ASP	N-CA-C	5.13	118.69	112.23
1	C	109	ILE	N-CA-C	5.08	115.85	110.72
1	A	48	TYR	N-CA-C	5.03	117.68	109.59
1	E	48	TYR	N-CA-C	5.02	117.93	109.95
1	A	132	VAL	N-CA-C	-5.01	101.09	108.11

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2440	0	2347	46	0
1	C	2440	0	2347	43	0
1	E	2440	0	2347	38	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1	0	0	0	0
3	A	19	0	13	5	0
3	C	19	0	13	4	0
3	E	19	0	13	2	0
4	A	183	0	0	2	0
4	C	176	0	0	1	0
4	E	185	0	0	1	0
All	All	7924	0	7080	126	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 9.

All (126) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:268:THR:HG21	3:C:311:BPX:HAD	1.52	0.91
1:A:268:THR:HG21	3:A:311:BPX:HAD	1.53	0.88
1:A:203:LEU:HD13	1:A:247:ILE:HD11	1.60	0.83
1:E:268:THR:HG21	3:E:311:BPX:HAD	1.63	0.79
1:A:243:ILE:O	1:A:247:ILE:HD12	1.83	0.79
1:A:133:THR:HG22	1:A:134:SER:N	1.98	0.78
1:C:133:THR:HG23	1:C:134:SER:N	1.97	0.78
1:C:205:PRO:HD3	1:C:266:SER:O	1.91	0.71
1:E:198:TYR:HA	1:E:271:LEU:O	1.89	0.71
1:E:204:TYR:HB2	1:E:205:PRO:HD2	1.75	0.68
1:E:224:LYS:O	1:E:228:GLU:HG3	1.95	0.66
1:E:243:ILE:HD11	1:E:255:ILE:HD11	1.76	0.66
1:A:204:TYR:HB2	1:A:205:PRO:HD2	1.79	0.65
1:C:204:TYR:HB2	1:C:205:PRO:HD2	1.79	0.64
1:C:105:ILE:HD13	1:C:301:MET:HE1	1.80	0.64
1:C:198:TYR:HA	1:C:271:LEU:O	1.99	0.63
1:A:168:LYS:HG2	1:A:169:TYR:HD2	1.63	0.63
1:A:198:TYR:HA	1:A:271:LEU:O	1.99	0.61
1:A:62:ILE:HD13	1:A:301:MET:HG2	1.82	0.61
1:A:119:THR:HA	1:A:123:ASN:O	2.00	0.60
1:E:62:ILE:HD13	1:E:301:MET:HG2	1.83	0.60
1:A:203:LEU:CD1	1:A:247:ILE:HD11	2.31	0.60
1:C:205:PRO:HB2	1:C:213:ILE:HG21	1.83	0.59
1:C:228:GLU:HG3	4:C:461:HOH:O	2.03	0.59
1:C:103:MET:HE1	1:C:304:THR:HG22	1.84	0.59
1:A:205:PRO:HB2	1:A:213:ILE:HG21	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:205:PRO:HD3	1:E:266:SER:O	2.04	0.58
1:A:243:ILE:HD11	1:A:255:ILE:HD11	1.85	0.57
1:E:71:ARG:HD3	1:E:125:LEU:O	2.05	0.56
1:E:288:PRO:O	1:E:292:GLU:HG2	2.06	0.56
1:A:88:GLU:O	1:A:92:GLN:HG3	2.05	0.56
1:A:224:LYS:O	1:A:227:VAL:HG22	2.06	0.56
1:C:74:ILE:HD13	1:C:281:LEU:HD23	1.87	0.56
1:A:224:LYS:O	1:A:228:GLU:HG3	2.06	0.55
1:A:42:TYR:OH	1:A:132:VAL:HG12	2.08	0.54
1:E:298:LEU:O	1:E:302:GLU:HG3	2.08	0.53
1:C:247:ILE:HG22	1:C:248:TYR:CG	2.44	0.53
1:C:272:ARG:O	1:C:273:ASP:HB3	2.08	0.53
1:A:205:PRO:HD3	1:A:266:SER:O	2.08	0.53
1:A:134:SER:OG	1:A:135:SER:N	2.41	0.53
1:C:247:ILE:HG22	1:C:248:TYR:CD2	2.43	0.52
1:A:74:ILE:HD13	1:A:281:LEU:HD23	1.91	0.52
1:A:272:ARG:O	1:A:273:ASP:HB3	2.09	0.52
1:C:74:ILE:CG2	1:C:197:SER:HB2	2.39	0.52
1:A:121:SER:HB2	1:C:247:ILE:O	2.10	0.52
1:C:270:GLU:OE2	3:C:311:BPX:HAF	2.11	0.51
1:C:103:MET:HE1	1:C:304:THR:CG2	2.40	0.51
1:C:132:VAL:HG12	1:C:132:VAL:O	2.09	0.51
1:C:62:ILE:HD13	1:C:301:MET:HE2	1.92	0.51
1:E:272:ARG:O	1:E:273:ASP:HB3	2.11	0.51
1:E:276:ARG:HD2	1:E:277:TYR:CE2	2.46	0.51
1:C:287:ILE:O	1:C:291:GLN:HG3	2.11	0.51
1:A:231:LYS:NZ	1:A:231:LYS:HB2	2.26	0.50
1:E:150:GLY:O	1:E:251:SER:HB2	2.12	0.50
1:C:119:THR:HA	1:C:123:ASN:O	2.12	0.50
1:A:273:ASP:HB2	1:A:281:LEU:HD13	1.93	0.50
1:E:270:GLU:OE2	3:E:311:BPX:HAL	2.11	0.50
1:E:198:TYR:O	1:E:199:SER:HB3	2.11	0.49
1:E:287:ILE:O	1:E:291:GLN:HG3	2.11	0.49
1:E:119:THR:HA	1:E:123:ASN:O	2.12	0.49
1:A:71:ARG:HD3	1:A:125:LEU:O	2.13	0.49
1:A:287:ILE:O	1:A:291:GLN:HG3	2.12	0.49
1:A:298:LEU:O	1:A:302:GLU:HG3	2.11	0.49
1:C:198:TYR:O	1:C:199:SER:HB3	2.14	0.48
1:A:270:GLU:OE2	3:A:311:BPX:HAF	2.13	0.48
1:E:160:PRO:HA	1:E:165:TYR:CD2	2.49	0.48
1:A:198:TYR:O	1:A:199:SER:HB3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:215:ASP:OD1	1:E:264:LYS:HE3	2.13	0.48
1:C:268:THR:HG21	3:C:311:BPX:CAD	2.34	0.47
1:A:231:LYS:HB2	1:A:231:LYS:HZ2	1.79	0.47
1:C:71:ARG:HD3	1:C:125:LEU:O	2.15	0.47
1:E:63:TRP:HE1	1:E:65:ASP:HB3	1.79	0.47
1:A:139:VAL:HG22	1:A:140:GLY:N	2.29	0.47
1:C:270:GLU:OE2	3:C:311:BPX:HAL	2.15	0.47
1:E:247:ILE:HG22	1:E:248:TYR:CG	2.51	0.46
1:C:204:TYR:HB2	1:C:205:PRO:CD	2.45	0.46
1:A:270:GLU:OE2	3:A:311:BPX:HAL	2.16	0.46
1:E:205:PRO:HB2	1:E:213:ILE:HG21	1.98	0.45
1:E:230:LEU:C	1:E:230:LEU:HD23	2.40	0.45
1:C:133:THR:HG23	1:C:134:SER:H	1.78	0.45
1:C:198:TYR:O	1:C:199:SER:CB	2.64	0.45
1:C:56:GLY:HA3	1:C:59:ARG:NE	2.31	0.45
3:A:311:BPX:OAN	3:A:311:BPX:HAJ	2.16	0.45
1:A:272:ARG:HH11	1:A:285:GLN:HE21	1.65	0.45
1:C:273:ASP:HB2	1:C:281:LEU:HD13	1.98	0.45
1:A:136:SER:HA	4:A:578:HOH:O	2.17	0.45
1:C:171:ASN:HB3	1:C:177:LYS:HD3	1.98	0.45
1:E:145:ARG:HG3	4:E:511:HOH:O	2.18	0.44
1:E:204:TYR:HB2	1:E:205:PRO:CD	2.44	0.44
1:A:204:TYR:HB2	1:A:205:PRO:CD	2.47	0.43
1:C:160:PRO:HA	1:C:165:TYR:CD2	2.53	0.43
1:C:297:VAL:O	1:C:301:MET:HG3	2.18	0.43
1:E:198:TYR:O	1:E:199:SER:CB	2.66	0.43
1:A:74:ILE:CG2	1:A:197:SER:HB2	2.48	0.43
1:A:302:GLU:O	1:A:305:VAL:HG12	2.18	0.43
1:E:180:VAL:HG13	1:E:257:TRP:CZ2	2.53	0.43
1:A:59:ARG:O	1:A:60:PRO:C	2.60	0.43
1:E:170:ALA:O	1:E:171:ASN:HB2	2.19	0.43
1:E:248:TYR:CD1	1:E:248:TYR:C	2.96	0.43
1:A:198:TYR:O	1:A:199:SER:CB	2.66	0.43
1:C:7:PHE:CE2	1:C:80:VAL:HG11	2.54	0.42
1:C:14:THR:OG1	1:C:17:GLU:HG3	2.19	0.42
1:E:302:GLU:O	1:E:305:VAL:HG12	2.20	0.42
1:C:298:LEU:O	1:C:302:GLU:HG3	2.19	0.42
1:E:281:LEU:HA	1:E:282:PRO:HD3	1.95	0.42
1:E:231:LYS:NZ	1:E:231:LYS:HB2	2.34	0.42
1:A:7:PHE:CE2	1:A:80:VAL:HG11	2.55	0.41
1:A:230:LEU:HD23	1:A:230:LEU:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:248:TYR:C	1:E:248:TYR:HD1	2.27	0.41
1:A:270:GLU:OE2	3:A:311:BPX:OAP	2.39	0.41
1:C:224:LYS:O	1:C:227:VAL:HG22	2.20	0.41
1:A:248:TYR:CD1	1:A:248:TYR:C	2.99	0.41
1:E:238:TYR:OH	1:E:271:LEU:HA	2.21	0.41
1:A:287:ILE:HB	1:A:288:PRO:HD3	2.03	0.41
1:C:105:ILE:CD1	1:C:301:MET:HE1	2.47	0.41
1:C:139:VAL:HG12	1:C:140:GLY:N	2.36	0.41
1:C:302:GLU:O	1:C:305:VAL:HG12	2.20	0.41
1:E:63:TRP:NE1	1:E:65:ASP:HB3	2.35	0.41
1:A:121:SER:CB	1:C:247:ILE:O	2.69	0.41
1:C:150:GLY:O	1:C:251:SER:HB2	2.21	0.41
1:A:303:HIS:HD2	4:A:454:HOH:O	2.04	0.40
1:E:171:ASN:HB3	1:E:177:LYS:HD3	2.03	0.40
1:A:5:ASN:HD22	1:A:5:ASN:HA	1.71	0.40
1:E:273:ASP:OD1	1:E:275:GLY:N	2.45	0.40
1:E:287:ILE:HB	1:E:288:PRO:HD3	2.03	0.40
1:C:85:LYS:HE2	1:C:294:TRP:CD1	2.57	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	305/307 (99%)	289 (95%)	12 (4%)	4 (1%)	9	2
1	C	305/307 (99%)	289 (95%)	12 (4%)	4 (1%)	9	2
1	E	305/307 (99%)	290 (95%)	10 (3%)	5 (2%)	7	2
All	All	915/921 (99%)	868 (95%)	34 (4%)	13 (1%)	9	2

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	SER
1	C	199	SER
1	A	205	PRO
1	C	205	PRO
1	E	198	TYR
1	E	199	SER
1	E	205	PRO
1	E	272	ARG
1	A	272	ARG
1	A	198	TYR
1	C	198	TYR
1	C	272	ARG
1	E	134	SER

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	264/264 (100%)	258 (98%)	6 (2%)	44	30
1	C	264/264 (100%)	256 (97%)	8 (3%)	36	21
1	E	264/264 (100%)	259 (98%)	5 (2%)	50	38
All	All	792/792 (100%)	773 (98%)	19 (2%)	43	28

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ASN
1	A	198	TYR
1	A	205	PRO
1	A	231	LYS
1	A	247	ILE
1	A	248	TYR
1	C	132	VAL
1	C	133	THR
1	C	134	SER
1	C	137	LEU

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Mol	Chain	Res	Type
1	C	198	TYR
1	C	205	PRO
1	C	228	GLU
1	C	248	TYR
1	E	2	ARG
1	E	198	TYR
1	E	205	PRO
1	E	231	LYS
1	E	248	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	37	GLN
1	A	92	GLN
1	A	112	ASN
1	A	221	GLN
1	A	285	GLN
1	A	303	HIS
1	C	5	ASN
1	C	37	GLN
1	C	92	GLN
1	C	112	ASN
1	C	221	GLN
1	C	261	GLN
1	C	285	GLN
1	E	5	ASN
1	E	37	GLN
1	E	58	ASN
1	E	92	GLN
1	E	112	ASN
1	E	261	GLN
1	E	285	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	BPX	E	311	2	16,19,19	2.38	1 (6%)	15,26,26	0.88	1 (6%)
3	BPX	C	311	2	16,19,19	2.39	1 (6%)	15,26,26	0.86	1 (6%)
3	BPX	A	311	2	16,19,19	2.40	1 (6%)	15,26,26	0.85	1 (6%)

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	BPX	E	311	2	-	4/17/18/18	0/1/1/1
3	BPX	C	311	2	-	4/17/18/18	0/1/1/1
3	BPX	A	311	2	-	4/17/18/18	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	311	BPX	OAS-NAM	9.09	1.42	1.22
3	C	311	BPX	OAS-NAM	9.08	1.42	1.22
3	E	311	BPX	OAS-NAM	9.04	1.42	1.22

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	311	BPX	OAQ-CAC-CAB	-2.34	116.65	122.86
3	C	311	BPX	OAQ-CAC-CAB	-2.30	116.76	122.86
3	E	311	BPX	OAQ-CAC-CAB	-2.23	116.95	122.86

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	311	BPX	CAF-CAB-CAC-OAQ
3	A	311	BPX	CAF-CAB-CAC-OAN
3	E	311	BPX	CAF-CAB-CAC-OAQ
3	E	311	BPX	CAF-CAB-CAC-OAN
3	E	311	BPX	CAG-CAA-CAB-CAC
3	C	311	BPX	CAG-CAA-CAB-CAC
3	A	311	BPX	CAG-CAA-CAB-CAF
3	E	311	BPX	CAG-CAA-CAB-CAF
3	C	311	BPX	CAF-CAB-CAC-OAQ
3	C	311	BPX	CAF-CAB-CAC-OAN
3	A	311	BPX	CAG-CAA-CAB-CAC
3	C	311	BPX	CAG-CAA-CAB-CAF

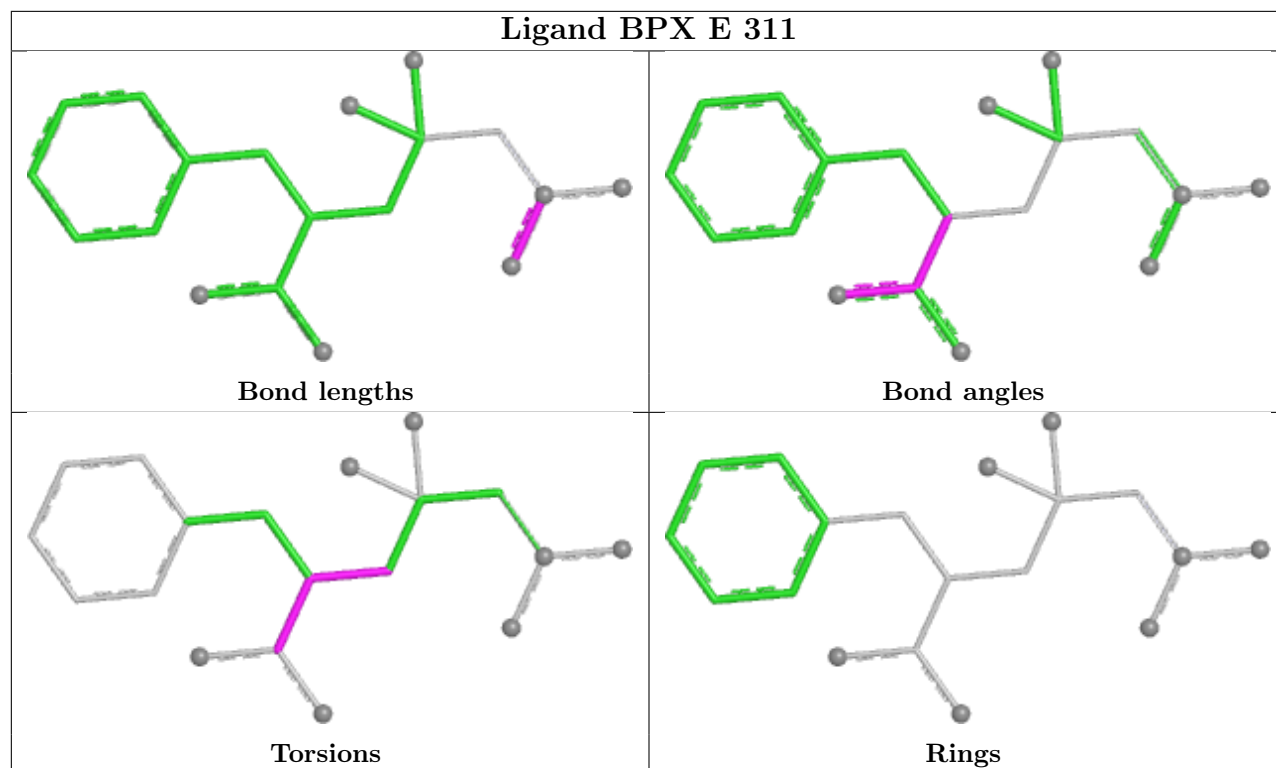
There are no ring outliers.

3 monomers are involved in 11 short contacts:

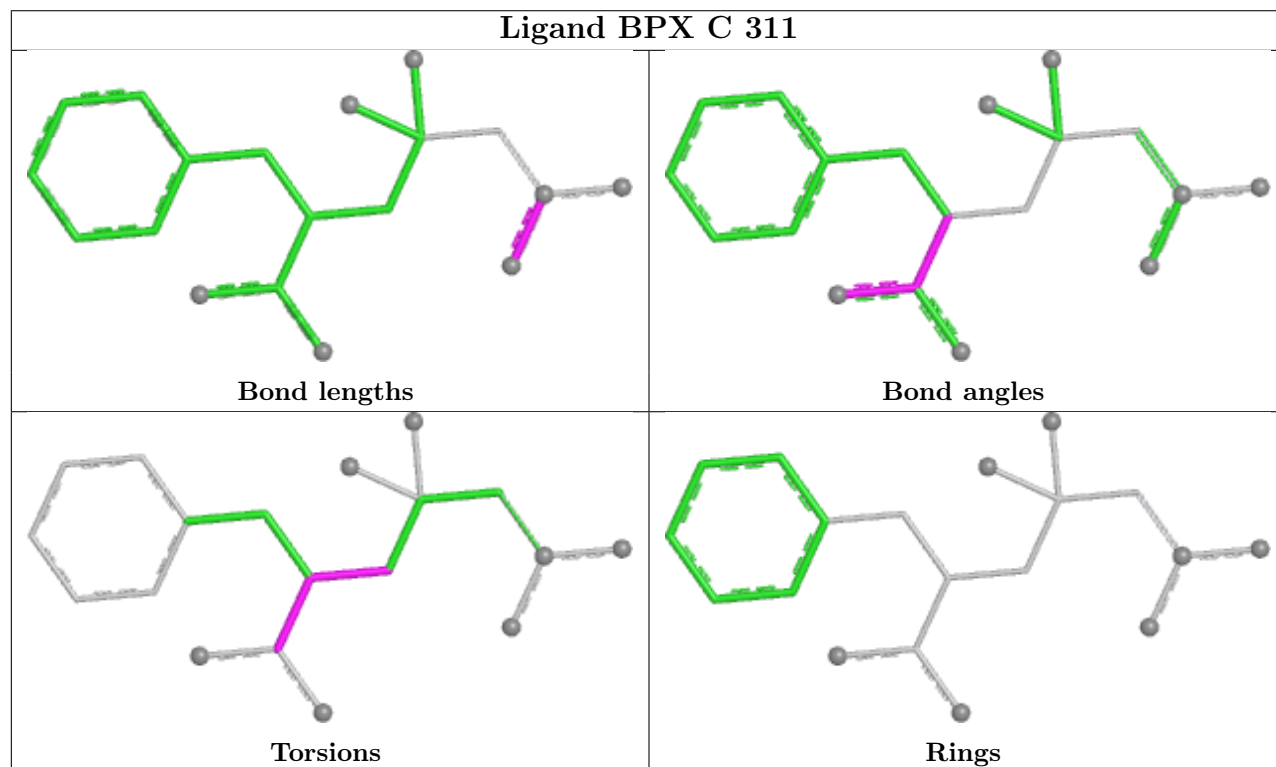
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	311	BPX	2	0
3	C	311	BPX	4	0
3	A	311	BPX	5	0

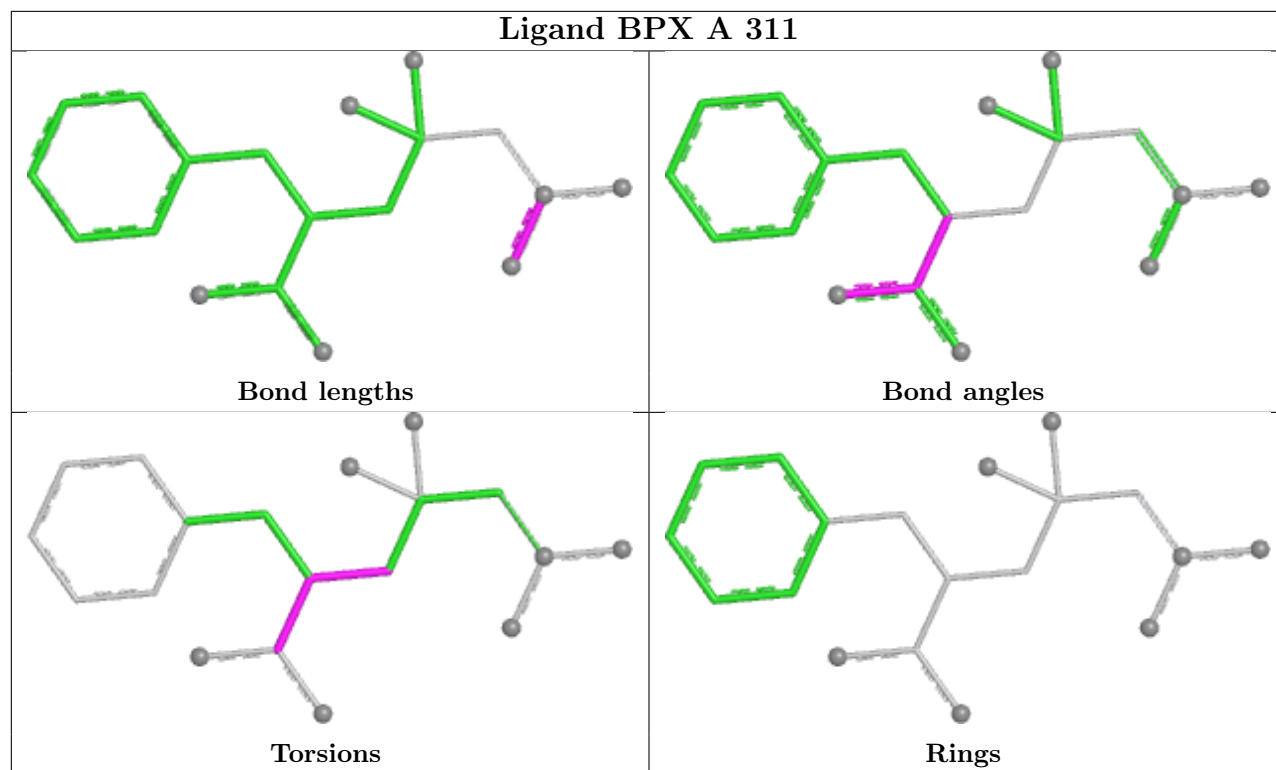
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand BPX E 311



Ligand BPX C 311





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	307/307 (100%)	0.47	26 (8%) 16 17	8, 14, 34, 82	3 (0%)
1	C	307/307 (100%)	0.78	36 (11%) 9 8	10, 16, 35, 96	3 (0%)
1	E	307/307 (100%)	0.67	28 (9%) 15 15	9, 16, 33, 84	3 (0%)
All	All	921/921 (100%)	0.64	90 (9%) 13 13	8, 15, 34, 96	9 (0%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1	ALA	7.6
1	A	307	ASN	6.7
1	A	137	LEU	6.5
1	A	135	SER	6.1
1	E	307	ASN	6.0
1	C	248	TYR	5.9
1	E	134	SER	5.7
1	E	205	PRO	5.7
1	C	133	THR	5.5
1	A	305	VAL	5.3
1	A	134	SER	5.2
1	A	247	ILE	5.1
1	C	247	ILE	5.1
1	C	305	VAL	5.0
1	C	132	VAL	5.0
1	C	205	PRO	4.9
1	C	307	ASN	4.9
1	C	135	SER	4.8
1	E	305	VAL	4.8
1	E	133	THR	4.7
1	C	1	ALA	4.7
1	A	2	ARG	4.7
1	E	137	LEU	4.6

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Mol	Chain	Res	Type	RSRZ
1	A	1	ALA	4.5
1	A	306	ASN	4.4
1	E	135	SER	4.4
1	A	205	PRO	4.3
1	A	248	TYR	4.3
1	C	137	LEU	4.2
1	C	2	ARG	4.0
1	E	197	SER	4.0
1	A	133	THR	3.9
1	E	136	SER	3.8
1	E	306	ASN	3.8
1	C	58	ASN	3.7
1	C	306	ASN	3.7
1	E	132	VAL	3.7
1	C	206	TYR	3.7
1	C	134	SER	3.5
1	C	197	SER	3.5
1	A	57	SER	3.4
1	A	136	SER	3.4
1	A	197	SER	3.4
1	E	276	ARG	3.4
1	E	57	SER	3.4
1	E	248	TYR	3.3
1	E	2	ARG	3.2
1	C	303	HIS	3.0
1	C	276	ARG	2.9
1	A	206	TYR	2.9
1	C	168	LYS	2.9
1	A	273	ASP	2.9
1	C	57	SER	2.8
1	E	199	SER	2.8
1	E	273	ASP	2.8
1	A	276	ARG	2.8
1	E	198	TYR	2.8
1	A	198	TYR	2.7
1	C	207	GLY	2.7
1	C	100	LEU	2.7
1	C	199	SER	2.7
1	E	228	GLU	2.6
1	E	58	ASN	2.6
1	E	303	HIS	2.5
1	A	138	CYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	304	THR	2.5
1	C	32	LEU	2.5
1	C	273	ASP	2.4
1	C	56	GLY	2.4
1	C	102	SER	2.4
1	A	132	VAL	2.4
1	C	55	GLY	2.4
1	C	91	GLY	2.4
1	A	199	SER	2.4
1	C	198	TYR	2.3
1	E	207	GLY	2.3
1	A	3	SER	2.3
1	E	138	CYS	2.3
1	C	60	PRO	2.3
1	A	92	GLN	2.2
1	C	261	GLN	2.2
1	E	56	GLY	2.2
1	A	58	ASN	2.2
1	E	247	ILE	2.1
1	C	225	SER	2.1
1	E	5	ASN	2.1
1	A	31	GLN	2.1
1	C	136	SER	2.0
1	E	206	TYR	2.0
1	C	155	GLY	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](#)

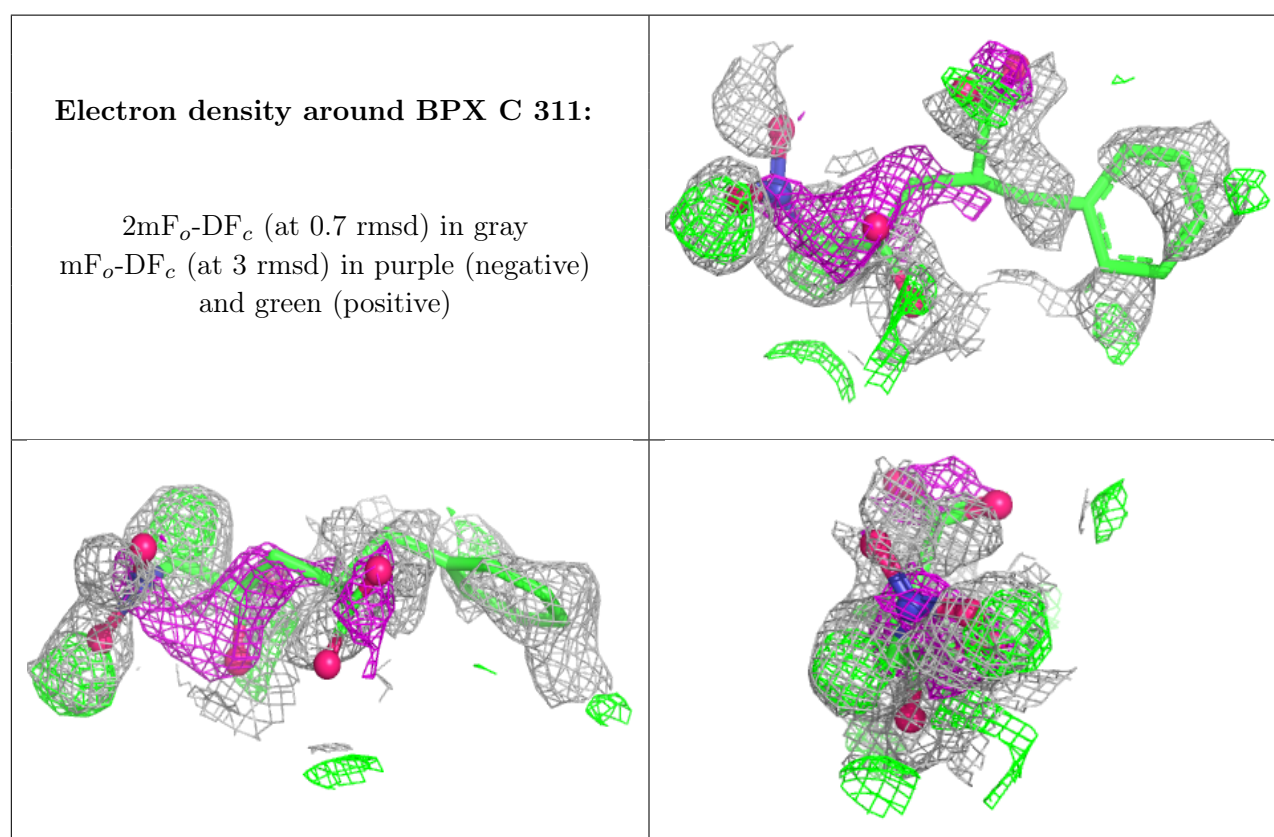
There are no oligosaccharides in this entry.

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.

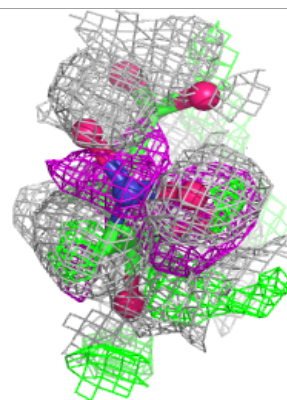
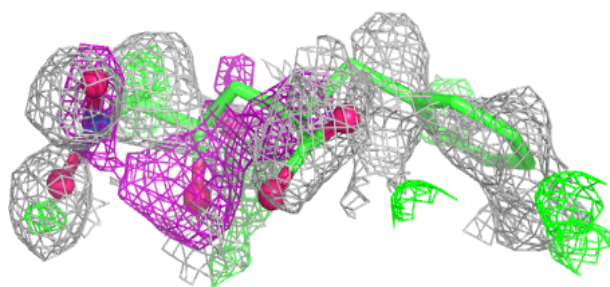
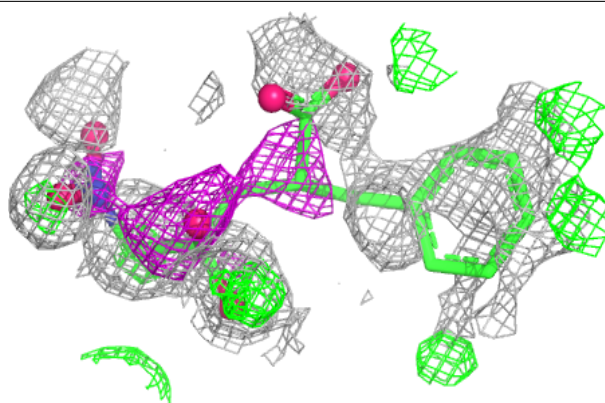
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BPX	C	311	19/19	0.56	0.25	44,49,51,51	0
3	BPX	E	311	19/19	0.64	0.23	39,50,51,51	0
3	BPX	A	311	19/19	0.67	0.24	39,48,49,49	0
2	ZN	E	309	1/1	0.98	0.04	14,14,14,14	0
2	ZN	C	309	1/1	0.99	0.04	14,14,14,14	0
2	ZN	A	309	1/1	1.00	0.02	11,11,11,11	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.

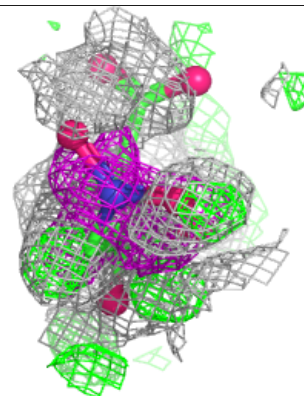
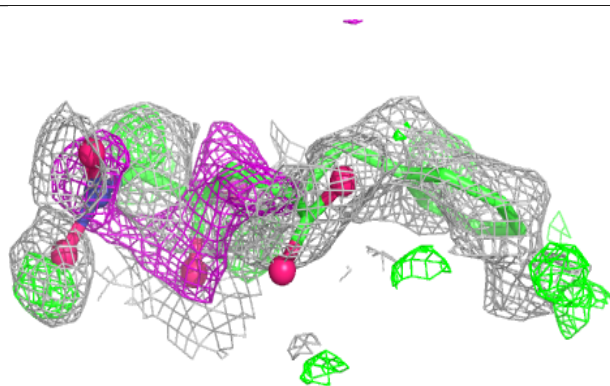
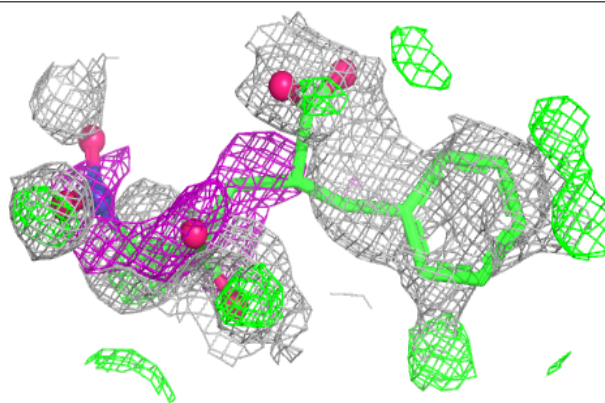


Electron density around BPX E 311:

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

**Electron density around BPX A 311:**

$2mF_o-DF_c$ (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.