



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:23 AM UTC

PDB ID : 2FZC / pdb_00002fzc
Title : The Structure of Wild-Type E. Coli Aspartate Transcarbamoylase in Complex with Novel T State Inhibitors at 2.10 Resolution
Authors : Heng, S.; Stieglitz, K.A.; Eldo, J.; Xia, J.; Cardia, J.P.; Kantrowitz, E.R.
Deposited on : 2006-02-09
Resolution : 2.10 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

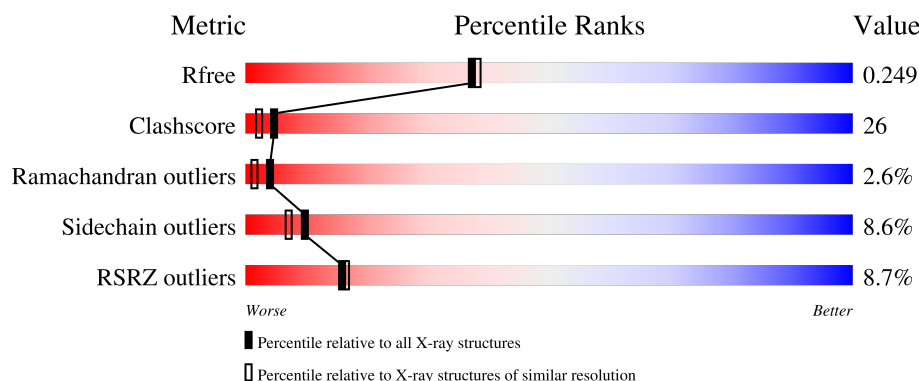
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	310	6% 67% 25% 6% .
1	C	310	5% 68% 27% . .
2	B	153	16% 52% 35% 10% . .
2	D	153	15% 48% 35% 13% . .

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	EOP	C	812	-	-	X	-
5	CTP	D	902	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8075 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.

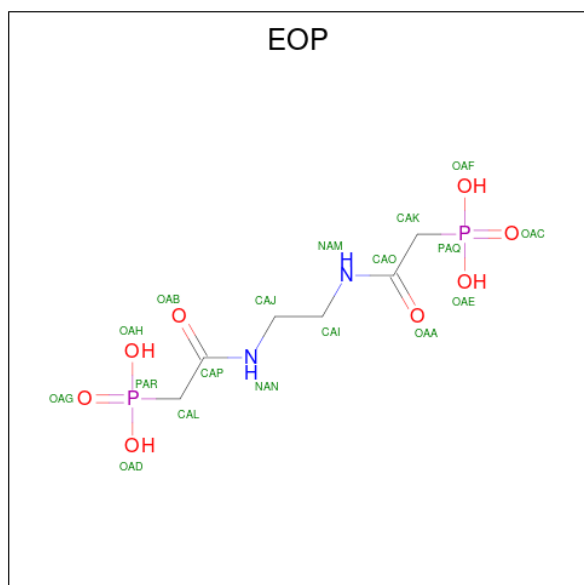
- Molecule 1 is a protein called Aspartate carbamoyltransferase catalytic chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			
1	C	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			

- Molecule 2 is a protein called Aspartate carbamoyltransferase regulatory chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	152	Total	C	N	O	S	0	0	0
			1191	746	212	228	5			
2	D	152	Total	C	N	O	S	0	0	0
			1191	746	212	228	5			

- Molecule 3 is {ETHANE-1,2-DIYLBIS[IMINO(2-OXOETHANE-2,1-DIYL)]}BIS(PHOSPHONIC ACID) (CCD ID: EOP) (formula: C₆H₁₄N₂O₈P₂).

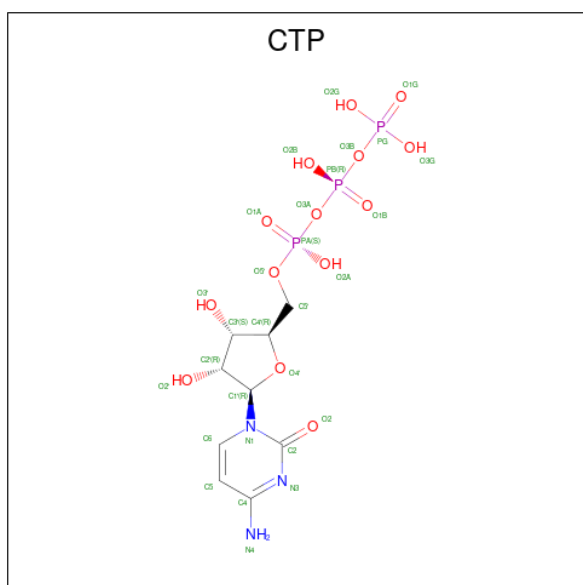


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			18	6	2	8	2		
3	C	1	Total	C	N	O	P	0	0
			18	6	2	8	2		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		

- Molecule 5 is CYTIDINE-5'-TRIPHOSPHATE (CCD ID: CTP) (formula: C₉H₁₆N₃O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			29	9	3	14	3		
5	D	1	Total	C	N	O	P	0	0
			29	9	3	14	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	234	Total	O	0	0
			234	234		

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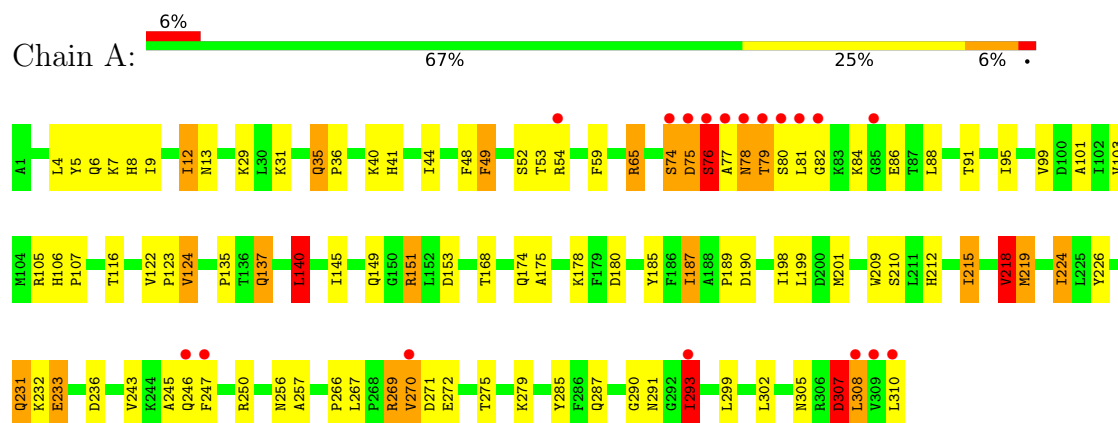
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	124	Total 124	O 124	0	0
6	C	284	Total 284	O 284	0	0
6	D	125	Total 125	O 125	0	0

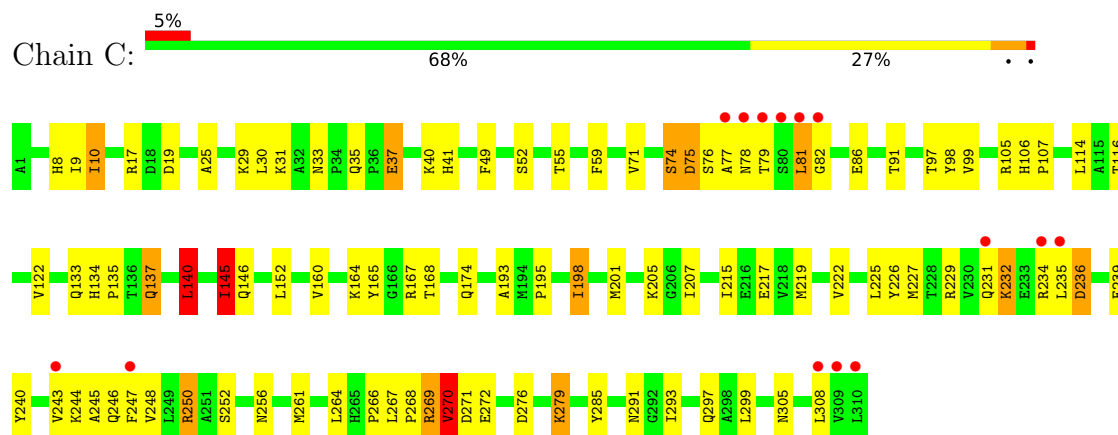
3 Residue-property plots [i](#)

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.

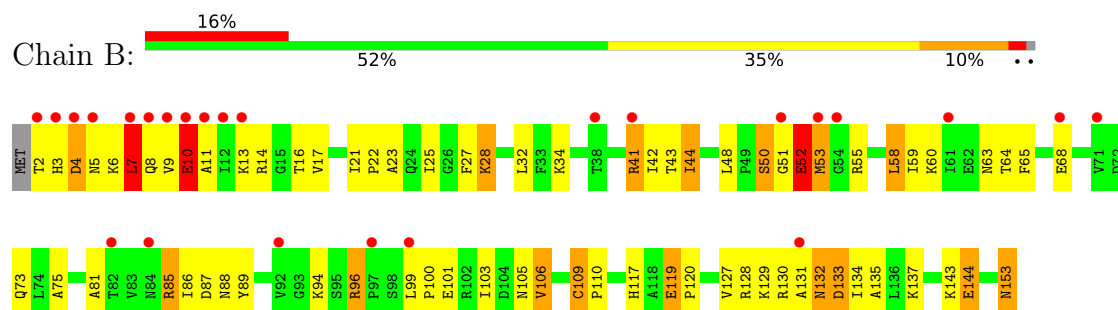
- Molecule 1: Aspartate carbamoyltransferase catalytic chain



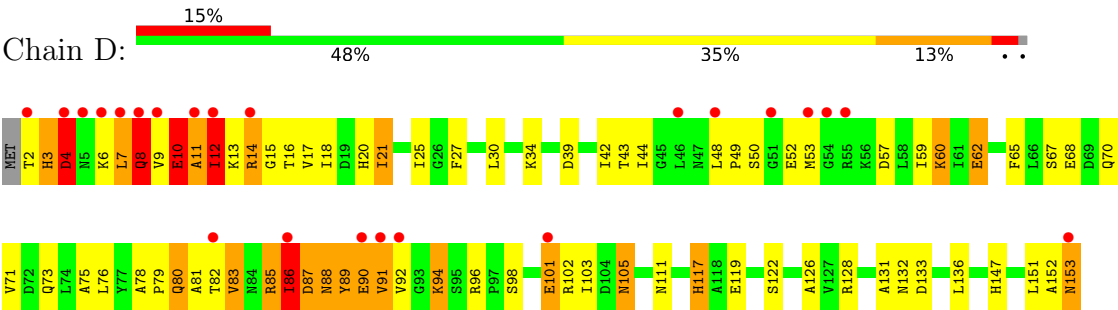
- Molecule 1: Aspartate carbamoyltransferase catalytic chain



- Molecule 2: Aspartate carbamoyltransferase regulatory chain



● Molecule 2: Aspartate carbamoyltransferase regulatory chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.59Å 120.59Å 141.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.35 – 2.10 27.35 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (27.35-2.10) 99.6 (27.35-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 1.99Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.250 0.201 , 0.249	Depositor DCC
R_{free} test set	8145 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	31.1	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 62.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8075	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% 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: CTP, ZN, EOP

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	0.86	3/2461 (0.1%)	1.25	22/3339 (0.7%)
1	C	0.93	0/2461	1.23	21/3339 (0.6%)
2	B	0.75	0/1209	1.26	13/1634 (0.8%)
2	D	0.72	0/1209	1.33	16/1634 (1.0%)
All	All	0.85	3/7340 (0.0%)	1.26	72/9946 (0.7%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	124	VAL	CA-CB	7.62	1.64	1.54
1	A	293	ILE	CA-CB	6.87	1.63	1.54
1	A	293	ILE	CG1-CD1	6.29	1.76	1.51

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	86	ILE	N-CA-C	15.05	140.64	109.34
1	C	77	ALA	N-CA-C	-12.54	96.47	112.90
2	B	119	GLU	N-CA-C	9.42	125.65	109.82
2	D	119	GLU	N-CA-C	9.21	124.19	110.10
1	A	269	ARG	N-CA-C	-8.75	96.21	109.94
1	C	122	VAL	N-CA-C	8.74	116.59	107.76
2	D	117	HIS	N-CA-C	8.10	119.74	111.07
2	D	90	GLU	N-CA-C	7.80	122.33	109.06
2	D	86	ILE	CA-C-N	7.35	135.57	121.54
2	D	86	ILE	C-N-CA	7.35	135.57	121.54
1	C	140	LEU	CA-CB-CG	7.22	141.56	116.30
1	C	269	ARG	N-CA-C	-7.16	98.76	109.86
2	B	10	GLU	N-CA-C	7.02	121.18	107.98
2	B	21	ILE	CA-C-N	-6.93	112.35	120.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	21	ILE	C-N-CA	-6.93	112.35	120.11
2	D	87	ASP	N-CA-C	6.92	125.53	110.80
2	B	117	HIS	N-CA-C	6.86	118.64	111.03
1	C	135	PRO	N-CA-C	6.80	123.67	113.81
1	A	307	ASP	CA-C-N	-6.80	112.86	123.13
1	A	307	ASP	C-N-CA	-6.80	112.86	123.13
1	A	257	ALA	N-CA-C	6.77	119.79	110.24
2	B	52	GLU	N-CA-C	-6.76	105.15	113.19
2	D	12	ILE	N-CA-C	6.63	123.13	109.34
2	D	83	VAL	N-CA-C	-6.54	99.01	108.36
2	B	75	ALA	N-CA-C	6.35	118.20	111.28
1	C	222	VAL	N-CA-C	6.27	117.67	109.58
2	B	44	ILE	N-CA-C	6.21	117.07	108.12
1	C	137	GLN	N-CA-C	-6.18	104.46	111.07
1	A	245	ALA	N-CA-C	6.15	118.77	111.33
1	A	140	LEU	CA-CB-CG	6.14	137.81	116.30
2	D	98	SER	N-CA-C	-6.05	101.04	110.42
1	A	99	VAL	N-CA-C	6.03	117.53	108.96
2	B	109	CYS	CA-C-N	6.00	125.62	119.56
2	B	109	CYS	C-N-CA	6.00	125.62	119.56
2	B	21	ILE	N-CA-C	-5.92	103.54	108.63
1	C	10	ILE	N-CA-C	5.85	116.64	111.90
1	A	53	THR	N-CA-C	5.81	118.09	111.11
1	A	137	GLN	N-CA-C	-5.79	105.05	111.36
1	A	180	ASP	N-CA-C	5.76	119.10	109.95
1	A	287	GLN	N-CA-C	-5.70	105.15	111.36
1	C	236	ASP	N-CA-C	-5.70	102.89	110.07
2	D	25	ILE	N-CA-C	-5.65	108.34	113.71
2	D	105	ASN	N-CA-C	5.65	122.83	110.80
1	A	178	LYS	N-CA-C	-5.64	106.40	113.28
1	C	133	GLN	N-CA-C	5.60	118.38	109.81
2	D	126	ALA	N-CA-C	-5.57	100.61	109.96
1	C	234	ARG	N-CA-C	-5.56	106.66	113.50
1	A	103	VAL	N-CA-C	-5.46	100.52	108.17
2	D	21	ILE	CB-CA-C	-5.46	105.86	110.94
1	A	135	PRO	N-CA-C	5.46	121.72	113.81
1	A	215	ILE	CB-CA-C	-5.46	104.78	112.14
1	C	217	GLU	CA-C-N	-5.45	116.04	122.36
1	C	217	GLU	C-N-CA	-5.45	116.04	122.36
1	C	49	PHE	N-CA-C	-5.38	105.52	111.71
1	C	145	ILE	CB-CA-C	-5.34	105.13	111.97
1	C	98	TYR	CB-CA-C	-5.34	100.75	109.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	ILE	CB-CG1-CD1	5.33	125.00	113.80
1	C	97	THR	N-CA-C	-5.33	106.78	113.28
2	D	62	GLU	N-CA-C	5.30	118.74	110.32
2	B	105	ASN	N-CA-C	5.28	122.05	110.80
1	A	49	PHE	N-CA-C	-5.25	105.67	111.71
2	D	75	ALA	N-CA-C	5.25	117.68	111.33
1	A	218	VAL	N-CA-C	5.24	119.55	113.42
1	A	74	SER	N-CA-C	5.24	114.95	108.19
1	C	71	VAL	N-CA-C	-5.23	100.34	108.81
1	A	35	GLN	CA-C-N	5.18	124.89	119.82
1	A	35	GLN	C-N-CA	5.18	124.89	119.82
1	C	19	ASP	N-CA-C	-5.12	105.59	111.07
1	C	99	VAL	N-CA-C	5.09	116.19	108.96
2	B	106	VAL	N-CA-C	5.04	119.46	112.35
1	A	88	LEU	N-CA-C	-5.02	105.81	111.28
1	C	193	ALA	N-CA-C	5.00	117.15	110.35

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	2415	0	2422	94	0
1	C	2415	0	2422	83	0
2	B	1191	0	1202	80	0
2	D	1191	0	1202	118	0
3	A	18	0	10	5	0
3	C	18	0	10	8	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	B	29	0	12	4	0
5	D	29	0	12	9	0
6	A	234	0	0	10	0
6	B	124	0	0	12	0
6	C	284	0	0	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	125	0	0	17	0
All	All	8075	0	7292	376	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 26.

All (376) 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:293:ILE:CG1	1:A:293:ILE:CD1	1.76	1.59
2:D:11:ALA:HB1	5:D:902:CTP:N3	1.36	1.37
2:B:96:ARG:NH1	2:B:96:ARG:HB3	1.61	1.16
1:C:33:ASN:HB2	6:C:873:HOH:O	1.50	1.11
2:D:101:GLU:CD	2:D:101:GLU:H	1.61	1.07
6:B:918:HOH:O	2:D:39:ASP:HB3	1.52	1.07
2:B:2:THR:HA	5:B:901:CTP:O2'	1.55	1.04
1:A:29:LYS:HD3	1:A:310:LEU:HB3	1.41	1.02
2:D:17:VAL:HG22	2:D:60:LYS:HG2	1.39	1.02
2:B:96:ARG:HB3	2:B:96:ARG:HH11	0.81	0.98
2:B:96:ARG:HH11	2:B:96:ARG:CB	1.75	0.97
1:C:10:ILE:HD11	1:C:116:THR:HG21	1.47	0.96
1:A:151:ARG:HG3	1:A:151:ARG:HH11	1.30	0.95
2:D:11:ALA:CB	5:D:902:CTP:N3	2.31	0.94
1:C:146:GLN:HG3	6:C:913:HOH:O	1.66	0.94
2:D:14:ARG:HB3	2:D:14:ARG:HH11	1.33	0.94
1:C:270:VAL:HG13	1:C:271:ASP:H	1.36	0.91
2:D:67:SER:H	2:D:70:GLN:HE21	1.20	0.90
6:B:952:HOH:O	2:D:10:GLU:HG2	1.71	0.89
2:D:4:ASP:HB3	2:D:8:GLN:HG2	1.53	0.88
2:D:67:SER:OG	2:D:70:GLN:HG3	1.76	0.85
2:D:86:ILE:HG23	2:D:87:ASP:N	1.91	0.85
2:B:4:ASP:OD1	2:B:9:VAL:HG21	1.77	0.85
1:C:75:ASP:HB3	6:C:955:HOH:O	1.77	0.84
1:A:81:LEU:HD12	1:A:86:GLU:OE2	1.78	0.84
5:D:902:CTP:H5'2	6:D:903:HOH:O	1.76	0.83
2:D:14:ARG:HH11	2:D:14:ARG:CB	1.96	0.78
2:B:27:PHE:HB3	2:B:28:LYS:HZ2	1.49	0.78
1:A:190:ASP:HB2	6:A:955:HOH:O	1.83	0.78
1:A:29:LYS:CD	1:A:310:LEU:HB3	2.14	0.78
2:D:9:VAL:HG13	2:D:43:THR:HG21	1.66	0.77
1:C:137:GLN:HE21	3:C:812:EOP:HAJ1	1.47	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:12:ILE:HG22	6:D:1017:HOH:O	1.84	0.77
2:D:85:ARG:HG2	2:D:86:ILE:N	1.99	0.76
2:D:86:ILE:HG12	2:D:87:ASP:H	1.49	0.76
1:A:52:SER:OG	3:A:811:EOP:HAL2	1.86	0.76
2:D:4:ASP:CB	2:D:8:GLN:HG2	2.15	0.76
2:B:22:PRO:HG2	2:B:25:ILE:HG13	1.66	0.76
2:B:153:ASN:HB2	6:B:902:HOH:O	1.85	0.75
1:C:81:LEU:HD22	1:C:91:THR:HG23	1.67	0.75
2:B:14:ARG:NH1	2:B:65:PHE:HE2	1.84	0.75
1:A:151:ARG:HG3	1:A:151:ARG:NH1	2.00	0.75
2:D:86:ILE:HG21	2:D:92:VAL:CG1	2.18	0.74
1:A:40:LYS:HG3	6:A:932:HOH:O	1.87	0.74
2:B:153:ASN:CB	6:B:902:HOH:O	2.36	0.74
1:A:81:LEU:HG	1:A:91:THR:HG23	1.67	0.74
2:B:81:ALA:O	2:B:96:ARG:HD2	1.88	0.74
2:D:101:GLU:CD	2:D:101:GLU:N	2.41	0.74
1:A:215:ILE:HG22	1:A:219:MET:HE2	1.68	0.73
1:C:105:ARG:NH2	3:C:812:EOP:HAL2	2.03	0.73
2:D:11:ALA:HB1	5:D:902:CTP:C2	2.23	0.72
1:C:270:VAL:HG13	1:C:271:ASP:N	2.04	0.72
5:D:902:CTP:O1A	6:D:903:HOH:O	2.09	0.71
2:D:85:ARG:HD3	2:D:86:ILE:HD12	1.73	0.71
1:C:305:ASN:HB2	1:C:308:LEU:HD21	1.73	0.71
2:B:13:LYS:HG3	2:B:14:ARG:H	1.54	0.70
1:C:10:ILE:HD11	1:C:116:THR:CG2	2.22	0.70
1:A:250:ARG:HG3	1:A:250:ARG:HH11	1.56	0.70
1:A:269:ARG:O	1:A:270:VAL:HB	1.90	0.70
5:D:902:CTP:PA	6:D:903:HOH:O	2.49	0.69
2:D:14:ARG:CB	2:D:14:ARG:NH1	2.55	0.69
2:B:133:ASP:OD2	2:B:133:ASP:C	2.35	0.69
2:D:4:ASP:HB3	2:D:8:GLN:CG	2.22	0.69
1:C:82:GLY:HA3	6:C:939:HOH:O	1.92	0.69
1:C:137:GLN:NE2	3:C:812:EOP:HAJ1	2.07	0.68
1:C:31:LYS:NZ	1:C:291:ASN:HD21	1.91	0.68
2:D:9:VAL:CG1	2:D:43:THR:HG21	2.23	0.68
2:D:12:ILE:HG12	2:D:62:GLU:OE1	1.92	0.68
1:A:302:LEU:CD2	1:A:308:LEU:HD22	2.23	0.68
2:B:3:HIS:O	2:B:9:VAL:HG13	1.94	0.68
2:D:86:ILE:CG2	2:D:92:VAL:HG12	2.25	0.67
2:B:153:ASN:ND2	6:B:1015:HOH:O	2.26	0.67
1:A:189:PRO:HG3	1:A:247:PHE:CE2	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:GLN:OE1	1:A:201:MET:HE2	1.96	0.66
2:B:99:LEU:HD22	2:B:127:VAL:HG11	1.76	0.66
2:B:101:GLU:HA	6:B:980:HOH:O	1.94	0.66
1:A:269:ARG:O	1:A:270:VAL:CB	2.43	0.66
2:D:60:LYS:HD2	6:D:913:HOH:O	1.94	0.66
1:C:246:GLN:HB2	1:C:247:PHE:CE2	2.30	0.66
2:D:52:GLU:HG2	2:D:53:MET:HG3	1.77	0.66
2:D:131:ALA:O	2:D:132:ASN:HB2	1.96	0.65
1:C:105:ARG:CZ	3:C:812:EOP:HAL2	2.27	0.65
2:D:67:SER:H	2:D:70:GLN:NE2	1.91	0.64
1:A:185:TYR:CD2	1:A:218:VAL:HG21	2.32	0.64
1:A:12:ILE:HD11	1:A:175:ALA:HB2	1.79	0.64
1:A:185:TYR:HD2	1:A:212:HIS:CE1	2.16	0.64
2:D:7:LEU:H	2:D:7:LEU:HD22	1.62	0.64
2:D:86:ILE:CG1	2:D:87:ASP:H	2.05	0.64
2:D:11:ALA:O	2:D:12:ILE:O	2.14	0.64
2:B:94:LYS:NZ	5:B:901:CTP:O1A	2.30	0.64
1:C:137:GLN:HA	1:C:140:LEU:HD13	1.79	0.64
1:C:279:LYS:HD2	1:C:279:LYS:N	2.12	0.64
1:A:231:GLN:HB3	1:A:233:GLU:OE2	1.97	0.63
2:D:27:PHE:HA	2:D:30:LEU:HD12	1.79	0.63
1:A:153:ASP:HB2	6:A:971:HOH:O	1.97	0.63
2:B:2:THR:OG1	2:B:9:VAL:CG1	2.46	0.63
2:D:82:THR:CB	2:D:96:ARG:NH1	2.62	0.62
1:A:75:ASP:O	1:A:77:ALA:N	2.31	0.62
1:A:151:ARG:HD2	6:A:891:HOH:O	1.98	0.62
2:B:68:GLU:CD	2:B:68:GLU:H	2.06	0.62
1:C:134:HIS:CD2	3:C:812:EOP:HAJ2	2.35	0.62
2:B:17:VAL:HG22	2:B:60:LYS:HG2	1.80	0.62
1:A:290:GLY:O	1:A:293:ILE:HG12	2.00	0.61
2:D:8:GLN:H	2:D:8:GLN:HE21	1.47	0.61
1:A:308:LEU:HD23	1:A:310:LEU:HD21	1.81	0.61
2:D:86:ILE:CG2	2:D:87:ASP:N	2.58	0.61
2:B:88:ASN:O	2:B:89:TYR:HB2	2.00	0.61
2:D:90:GLU:CD	2:D:91:VAL:H	2.08	0.61
2:B:51:GLY:C	2:B:53:MET:H	2.05	0.60
2:D:86:ILE:HG21	2:D:92:VAL:HG12	1.81	0.60
1:A:187:ILE:N	1:A:187:ILE:HD12	2.17	0.60
2:B:14:ARG:HH11	2:B:65:PHE:HE2	1.49	0.60
2:D:65:PHE:HD2	2:D:85:ARG:HG3	1.65	0.60
1:A:54:ARG:HG3	1:A:54:ARG:HH11	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ILE:HG23	1:A:224:ILE:HD11	1.83	0.60
1:A:308:LEU:HB3	1:A:310:LEU:HG	1.83	0.60
1:A:308:LEU:CG	1:A:310:LEU:HD21	2.31	0.60
6:B:1003:HOH:O	2:D:48:LEU:HD21	2.00	0.60
2:B:2:THR:CA	5:B:901:CTP:O2'	2.39	0.60
2:B:10:GLU:H	2:B:10:GLU:CD	2.08	0.60
1:C:31:LYS:HZ2	1:C:291:ASN:HD21	1.47	0.60
1:C:82:GLY:HA2	1:C:86:GLU:HB3	1.83	0.60
1:C:250:ARG:HB3	1:C:250:ARG:CZ	2.32	0.60
1:C:232:LYS:HE2	6:C:1045:HOH:O	2.00	0.60
1:A:308:LEU:HB3	1:A:310:LEU:CG	2.32	0.60
2:B:27:PHE:HB3	2:B:28:LYS:NZ	2.17	0.60
1:C:75:ASP:CB	6:C:955:HOH:O	2.43	0.59
1:C:270:VAL:CG1	6:C:921:HOH:O	2.50	0.59
1:A:145:ILE:HG23	1:A:224:ILE:CD1	2.31	0.59
1:C:9:ILE:HG21	1:C:299:LEU:HD21	1.84	0.59
1:C:246:GLN:O	1:C:247:PHE:CD1	2.55	0.59
1:C:10:ILE:CD1	1:C:116:THR:HG21	2.28	0.59
1:A:302:LEU:HD23	1:A:308:LEU:HD22	1.85	0.58
1:C:235:LEU:HD22	1:C:239:GLU:OE1	2.02	0.58
2:D:48:LEU:CD1	2:D:49:PRO:HD2	2.33	0.58
1:C:270:VAL:HG13	6:C:921:HOH:O	2.04	0.58
2:B:50:SER:OG	2:B:52:GLU:HG2	2.04	0.58
1:C:29:LYS:NZ	6:C:946:HOH:O	2.37	0.58
1:A:270:VAL:HG13	1:A:271:ASP:N	2.17	0.58
2:B:2:THR:OG1	2:B:9:VAL:HG12	2.04	0.57
2:B:130:ARG:NH1	2:B:133:ASP:OD2	2.37	0.57
2:B:41:ARG:NH1	2:B:41:ARG:HG3	2.19	0.57
1:C:245:ALA:HB2	6:C:993:HOH:O	2.04	0.57
1:C:17:ARG:NH1	6:C:902:HOH:O	2.36	0.57
1:A:308:LEU:HB3	1:A:310:LEU:HD11	1.86	0.57
1:C:114:LEU:C	1:C:114:LEU:HD23	2.30	0.57
2:D:90:GLU:OE2	2:D:90:GLU:HA	2.03	0.57
1:A:308:LEU:CD2	1:A:310:LEU:HD21	2.34	0.57
2:B:8:GLN:HB2	2:D:10:GLU:HA	1.87	0.57
2:B:41:ARG:HG3	2:B:41:ARG:HH11	1.69	0.57
2:D:60:LYS:HE3	6:D:913:HOH:O	2.04	0.57
2:D:117:HIS:HD2	6:D:981:HOH:O	1.87	0.57
1:A:8:HIS:C	1:A:9:ILE:HD12	2.30	0.57
1:A:308:LEU:O	1:A:310:LEU:HG	2.05	0.57
2:B:10:GLU:O	2:B:60:LYS:NZ	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ILE:HG21	1:A:299:LEU:HD21	1.86	0.57
2:D:86:ILE:HG22	2:D:90:GLU:O	2.05	0.57
1:C:146:GLN:CG	6:C:913:HOH:O	2.39	0.56
1:A:293:ILE:HD12	6:A:829:HOH:O	2.05	0.56
1:A:308:LEU:HG	1:A:310:LEU:HD21	1.87	0.56
1:A:250:ARG:HG3	1:A:250:ARG:NH1	2.21	0.56
1:C:55:THR:OG1	3:C:812:EOP:HAL1	2.05	0.56
1:A:31:LYS:NZ	1:A:291:ASN:HD21	2.04	0.56
2:D:14:ARG:HB3	2:D:14:ARG:NH1	2.11	0.55
1:C:37:GLU:HG3	1:C:40:LYS:HD3	1.89	0.55
1:C:74:SER:O	1:C:75:ASP:O	2.25	0.55
1:A:76:SER:C	1:A:78:ASN:N	2.65	0.55
2:B:32:LEU:HD22	2:B:106:VAL:HB	1.89	0.55
2:B:51:GLY:C	2:B:53:MET:N	2.64	0.55
2:D:6:LYS:HB3	2:D:7:LEU:HD22	1.89	0.55
1:A:308:LEU:HB3	1:A:310:LEU:CD1	2.37	0.55
1:A:137:GLN:HA	1:A:140:LEU:HD13	1.89	0.54
1:A:76:SER:C	1:A:78:ASN:H	2.14	0.54
2:D:14:ARG:HG2	2:D:65:PHE:HZ	1.73	0.54
3:A:811:EOP:OAC	3:A:811:EOP:OAD	2.24	0.54
1:C:270:VAL:CG1	1:C:271:ASP:H	2.16	0.54
2:D:12:ILE:CD1	6:D:936:HOH:O	2.55	0.54
2:D:49:PRO:HD2	6:D:935:HOH:O	2.07	0.54
1:C:215:ILE:HG22	1:C:219:MET:HE2	1.90	0.54
1:A:185:TYR:CG	1:A:218:VAL:HG21	2.43	0.53
1:C:168:THR:HB	6:C:852:HOH:O	2.08	0.53
1:A:305:ASN:HD22	1:A:308:LEU:HD12	1.73	0.53
2:B:2:THR:OG1	2:B:9:VAL:HG11	2.08	0.53
1:C:195:PRO:HD2	1:C:198:ILE:HD13	1.90	0.53
2:B:131:ALA:O	2:B:132:ASN:HB2	2.08	0.53
2:B:85:ARG:C	2:B:86:ILE:HD12	2.34	0.53
1:A:31:LYS:HZ1	1:A:291:ASN:HD21	1.57	0.52
2:D:34:LYS:NZ	6:D:1000:HOH:O	2.43	0.52
1:C:246:GLN:C	1:C:247:PHE:CG	2.87	0.52
2:B:2:THR:HG21	2:B:11:ALA:HB2	1.91	0.52
2:D:3:HIS:O	2:D:4:ASP:HB2	2.09	0.52
1:A:9:ILE:HG12	1:A:299:LEU:HD11	1.92	0.52
2:D:81:ALA:O	2:D:96:ARG:HD2	2.08	0.52
2:D:86:ILE:HD13	2:D:92:VAL:HG11	1.92	0.52
1:C:74:SER:C	1:C:75:ASP:O	2.53	0.51
1:A:74:SER:O	1:A:75:ASP:C	2.52	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:23:ALA:HB1	6:B:975:HOH:O	2.09	0.51
2:B:44:ILE:HB	2:D:44:ILE:HB	1.92	0.51
1:C:205:LYS:HE2	6:C:1089:HOH:O	2.10	0.51
2:D:7:LEU:HD22	2:D:7:LEU:N	2.25	0.51
1:A:290:GLY:O	1:A:293:ILE:CD1	2.59	0.51
2:B:4:ASP:CG	2:B:9:VAL:HG21	2.36	0.51
2:D:2:THR:HG23	2:D:50:SER:HA	1.92	0.51
2:D:48:LEU:HD12	6:D:935:HOH:O	2.11	0.51
1:A:215:ILE:CG2	1:A:219:MET:HE2	2.38	0.51
2:B:41:ARG:HD2	2:B:41:ARG:C	2.36	0.51
2:B:96:ARG:NH1	2:B:96:ARG:CB	2.52	0.51
1:A:305:ASN:ND2	1:A:307:ASP:O	2.44	0.51
2:D:14:ARG:NH1	2:D:14:ARG:HB2	2.25	0.51
2:D:85:ARG:CG	2:D:86:ILE:N	2.73	0.50
2:D:13:LYS:HG3	2:D:89:TYR:CE1	2.46	0.50
2:D:111:ASN:O	2:D:117:HIS:HE1	1.93	0.50
2:B:73:GLN:HE22	2:B:103:ILE:HD13	1.74	0.50
2:D:86:ILE:HB	2:D:92:VAL:HG12	1.92	0.50
1:A:49:PHE:O	1:A:76:SER:HB2	2.11	0.50
1:A:210:SER:OG	1:A:212:HIS:HE1	1.94	0.49
1:A:308:LEU:C	1:A:310:LEU:HG	2.36	0.49
2:B:8:GLN:HB2	2:D:10:GLU:CA	2.41	0.49
2:B:153:ASN:HB3	6:B:902:HOH:O	2.05	0.49
2:D:20:HIS:NE2	5:D:902:CTP:O1G	2.35	0.49
1:C:25:ALA:O	1:C:29:LYS:HG3	2.11	0.49
1:C:269:ARG:C	1:C:270:VAL:HG12	2.36	0.49
2:D:80:GLN:H	2:D:80:GLN:NE2	2.10	0.49
2:D:94:LYS:HG3	6:D:931:HOH:O	2.13	0.49
2:D:13:LYS:NZ	2:D:87:ASP:OD1	2.45	0.49
1:C:250:ARG:NE	1:C:276:ASP:OD2	2.38	0.49
2:D:73:GLN:HE22	2:D:103:ILE:HG23	1.78	0.49
2:D:8:GLN:HE21	2:D:8:GLN:N	2.11	0.49
2:D:80:GLN:H	2:D:80:GLN:CD	2.19	0.49
1:A:9:ILE:HD12	1:A:9:ILE:N	2.27	0.48
2:D:48:LEU:HD13	2:D:49:PRO:HD2	1.95	0.48
2:B:13:LYS:HG3	2:B:14:ARG:N	2.24	0.48
1:C:236:ASP:HB3	1:C:239:GLU:HB2	1.94	0.48
2:D:18:ILE:HG12	2:D:83:VAL:HG22	1.96	0.48
1:C:269:ARG:O	1:C:270:VAL:CB	2.61	0.48
2:D:60:LYS:CE	6:D:913:HOH:O	2.60	0.48
2:D:105:ASN:ND2	2:D:122:SER:OG	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:GLN:CB	1:A:233:GLU:OE2	2.62	0.47
1:A:266:PRO:HB2	3:A:811:EOP:HAI2	1.96	0.47
2:B:42:ILE:HG23	6:B:911:HOH:O	2.14	0.47
1:C:269:ARG:O	1:C:270:VAL:HB	2.14	0.47
2:D:60:LYS:CD	6:D:913:HOH:O	2.58	0.47
2:D:152:ALA:O	2:D:153:ASN:O	2.31	0.47
1:C:52:SER:OG	3:C:812:EOP:CAL	2.63	0.47
1:C:267:LEU:HB3	1:C:268:PRO:HA	1.95	0.47
1:C:40:LYS:O	1:C:41:HIS:HB2	2.15	0.47
1:A:40:LYS:O	1:A:41:HIS:HB2	2.14	0.47
1:C:271:ASP:HB2	6:C:921:HOH:O	2.13	0.47
2:B:41:ARG:HH11	2:B:41:ARG:CG	2.28	0.47
1:C:293:ILE:O	1:C:297:GLN:HG3	2.15	0.47
2:D:48:LEU:HD12	2:D:49:PRO:HD2	1.96	0.47
1:A:290:GLY:O	1:A:293:ILE:CG1	2.62	0.47
2:D:79:PRO:HG2	2:D:80:GLN:HE21	1.79	0.47
1:A:48:PHE:O	1:A:74:SER:HA	2.15	0.47
1:C:229:ARG:NH2	1:C:231:GLN:HG2	2.30	0.47
1:C:246:GLN:HB2	1:C:247:PHE:CZ	2.50	0.46
2:D:3:HIS:O	2:D:4:ASP:CB	2.63	0.46
1:A:149:GLN:HG3	1:A:224:ILE:HD11	1.97	0.46
2:B:14:ARG:HA	2:B:86:ILE:O	2.16	0.46
2:B:28:LYS:HE3	6:D:968:HOH:O	2.15	0.46
2:D:8:GLN:H	2:D:8:GLN:NE2	2.13	0.46
2:D:86:ILE:CB	2:D:92:VAL:HG12	2.46	0.46
1:A:270:VAL:CG1	1:A:271:ASP:N	2.78	0.46
1:C:174:GLN:OE1	1:C:201:MET:HE2	2.16	0.46
2:D:21:ILE:HB	2:D:57:ASP:O	2.16	0.46
1:A:91:THR:O	1:A:95:ILE:HG12	2.16	0.46
2:B:14:ARG:NH1	2:B:65:PHE:CE2	2.73	0.46
1:A:105:ARG:NH2	3:A:811:EOP:OAH	2.49	0.46
2:B:16:THR:OG1	2:B:65:PHE:HA	2.16	0.46
1:C:8:HIS:HB3	1:C:10:ILE:HD11	1.98	0.45
2:D:82:THR:CB	2:D:96:ARG:HH12	2.29	0.45
1:A:302:LEU:HD21	1:A:308:LEU:HD22	1.98	0.45
1:A:199:LEU:HD22	1:A:209:TRP:CE2	2.52	0.45
2:D:87:ASP:O	2:D:88:ASN:C	2.59	0.45
1:C:145:ILE:CD1	1:C:264:LEU:HD11	2.47	0.45
2:B:7:LEU:HB3	2:B:8:GLN:H	1.55	0.45
2:B:99:LEU:HA	2:B:99:LEU:HD23	1.60	0.45
1:A:12:ILE:HD11	1:A:175:ALA:CB	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:GLN:HG3	6:A:983:HOH:O	2.17	0.45
1:C:105:ARG:HH22	3:C:812:EOP:CAP	2.30	0.45
2:D:11:ALA:HB1	5:D:902:CTP:C4	2.36	0.45
1:A:81:LEU:HD21	1:A:95:ILE:HD11	1.98	0.44
1:A:81:LEU:HD11	1:A:91:THR:HA	1.99	0.44
1:C:219:MET:SD	1:C:261:MET:HE3	2.57	0.44
1:C:305:ASN:CB	1:C:308:LEU:HD21	2.45	0.44
2:D:65:PHE:CD2	2:D:85:ARG:HG3	2.49	0.44
2:B:16:THR:OG1	2:B:64:THR:O	2.33	0.44
2:B:73:GLN:NE2	6:B:1018:HOH:O	2.50	0.44
2:D:42:ILE:CG2	2:D:44:ILE:HD11	2.47	0.44
1:C:207:ILE:HD12	6:C:826:HOH:O	2.17	0.44
1:C:232:LYS:HB3	1:C:240:TYR:CE2	2.53	0.44
2:D:14:ARG:O	2:D:62:GLU:HB2	2.17	0.44
1:A:279:LYS:HE3	6:A:998:HOH:O	2.17	0.44
1:C:37:GLU:CG	1:C:40:LYS:HD3	2.47	0.44
1:C:79:THR:HG22	1:C:81:LEU:H	1.83	0.44
2:D:102:ARG:HD3	6:D:943:HOH:O	2.17	0.44
2:D:131:ALA:O	2:D:132:ASN:CB	2.61	0.44
1:A:77:ALA:HB3	6:A:1036:HOH:O	2.17	0.43
2:D:12:ILE:HG21	2:D:62:GLU:OE1	2.18	0.43
2:B:99:LEU:HD23	2:B:100:PRO:HD2	2.00	0.43
2:B:109:CYS:HA	2:B:110:PRO:HD3	1.81	0.43
2:D:128:ARG:HG3	6:D:1001:HOH:O	2.18	0.43
1:A:210:SER:OG	1:A:212:HIS:CE1	2.72	0.43
2:B:2:THR:CB	2:B:9:VAL:HG12	2.49	0.43
2:B:14:ARG:HB2	2:B:87:ASP:HA	2.00	0.43
2:D:15:GLY:HA3	2:D:62:GLU:HA	2.00	0.43
1:C:168:THR:CG2	1:C:226:TYR:OH	2.66	0.43
2:D:136:LEU:HD12	2:D:147:HIS:HA	2.00	0.43
2:B:63:ASN:N	2:B:63:ASN:ND2	2.66	0.43
2:D:39:ASP:OD1	2:D:39:ASP:N	2.42	0.43
2:D:44:ILE:CD1	2:D:59:ILE:HG23	2.49	0.43
1:C:248:VAL:HG22	1:C:271:ASP:O	2.18	0.43
1:C:256:ASN:OD1	1:C:256:ASN:N	2.52	0.43
1:A:36:PRO:HA	1:A:65:ARG:O	2.18	0.43
1:A:185:TYR:HD2	1:A:212:HIS:HE1	1.64	0.43
2:B:119:GLU:HB3	2:B:120:PRO:HD2	2.00	0.43
2:D:16:THR:CG2	2:D:83:VAL:HG13	2.49	0.43
2:D:94:LYS:HE2	2:D:94:LYS:O	2.19	0.43
1:A:13:ASN:HB3	6:A:1034:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:58:LEU:HD23	2:B:59:ILE:N	2.34	0.42
2:B:143:LYS:HD3	6:B:960:HOH:O	2.17	0.42
2:B:8:GLN:NE2	2:D:9:VAL:O	2.52	0.42
2:D:13:LYS:HB3	2:D:13:LYS:HE3	1.88	0.42
2:D:43:THR:C	2:D:44:ILE:HD13	2.43	0.42
2:D:152:ALA:O	2:D:153:ASN:C	2.62	0.42
2:B:88:ASN:O	2:B:89:TYR:CB	2.68	0.42
1:C:145:ILE:HD13	1:C:264:LEU:HD11	2.01	0.42
1:A:275:THR:O	1:A:279:LYS:NZ	2.46	0.42
2:D:73:GLN:NE2	2:D:103:ILE:HG23	2.35	0.42
1:A:4:LEU:HD23	1:A:7:LYS:HD3	2.01	0.42
1:C:30:LEU:CD2	1:C:35:GLN:HG3	2.49	0.42
2:D:78:ALA:HB1	2:D:81:ALA:HB2	2.02	0.42
2:B:58:LEU:HD21	2:B:60:LYS:HE3	2.02	0.42
2:B:130:ARG:HH11	2:B:130:ARG:HB3	1.85	0.42
1:C:75:ASP:OD1	1:C:76:SER:N	2.52	0.42
1:C:160:VAL:HG21	1:C:225:LEU:HD11	2.01	0.42
2:D:76:LEU:HD12	2:D:103:ILE:HD13	2.02	0.42
1:A:54:ARG:NH2	3:A:811:EOP:HAN	2.18	0.42
1:A:122:VAL:HA	1:A:123:PRO:HD3	1.91	0.42
2:B:48:LEU:O	2:B:55:ARG:HA	2.20	0.42
1:C:227:MET:O	1:C:266:PRO:HD2	2.19	0.42
1:A:232:LYS:HD3	1:A:232:LYS:C	2.44	0.42
1:C:250:ARG:NH1	1:C:252:SER:OG	2.52	0.42
2:B:133:ASP:OD2	2:B:134:ILE:N	2.52	0.41
1:C:164:LYS:HE2	1:C:165:TYR:CZ	2.55	0.41
1:A:5:TYR:O	1:A:6:GLN:HB2	2.20	0.41
5:B:901:CTP:O1B	5:B:901:CTP:H5'1	2.20	0.41
2:D:8:GLN:HB2	2:D:9:VAL:H	1.40	0.41
2:D:60:LYS:CB	2:D:60:LYS:NZ	2.84	0.41
1:A:44:ILE:HD12	1:A:101:ALA:HB3	2.02	0.41
2:D:2:THR:CG2	2:D:50:SER:HA	2.50	0.41
1:C:106:HIS:ND1	1:C:107:PRO:HD2	2.34	0.41
1:A:82:GLY:C	6:A:915:HOH:O	2.63	0.41
2:D:111:ASN:O	2:D:117:HIS:CE1	2.71	0.41
2:B:130:ARG:O	2:B:130:ARG:HG2	2.19	0.41
2:D:90:GLU:CD	2:D:91:VAL:N	2.77	0.41
2:D:90:GLU:CG	2:D:91:VAL:N	2.83	0.41
1:A:168:THR:CG2	1:A:226:TYR:OH	2.69	0.41
1:A:279:LYS:HE2	1:A:279:LYS:H	1.86	0.41
2:B:43:THR:HB	2:B:60:LYS:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:137:LYS:HB2	2:B:144:GLU:CG	2.51	0.41
1:C:152:LEU:HD23	1:C:152:LEU:HA	1.89	0.41
2:D:71:VAL:HG22	2:D:83:VAL:HG21	2.03	0.41
2:D:11:ALA:CB	5:D:902:CTP:C2	2.97	0.41
2:D:132:ASN:O	2:D:133:ASP:HB3	2.21	0.41
1:A:106:HIS:ND1	1:A:107:PRO:HD2	2.36	0.40
1:C:137:GLN:NE2	6:C:867:HOH:O	2.53	0.40
1:C:243:VAL:O	1:C:244:LYS:C	2.64	0.40
1:C:291:ASN:HD22	1:C:291:ASN:HA	1.73	0.40
1:A:76:SER:OG	1:A:79:THR:HB	2.20	0.40
2:B:86:ILE:HD12	2:B:86:ILE:N	2.37	0.40
2:B:130:ARG:O	2:B:131:ALA:C	2.63	0.40
2:B:129:LYS:HB2	2:B:129:LYS:HE2	1.75	0.40
2:B:128:ARG:HB3	2:B:135:ALA:HB3	2.04	0.40
1:A:266:PRO:O	1:A:267:LEU:HB2	2.21	0.40
2:D:10:GLU:H	2:D:10:GLU:HG3	1.59	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	308/310 (99%)	284 (92%)	17 (6%)	7 (2%)	5	2
1	C	308/310 (99%)	294 (96%)	12 (4%)	2 (1%)	21	18
2	B	150/153 (98%)	131 (87%)	14 (9%)	5 (3%)	3	1
2	D	150/153 (98%)	127 (85%)	13 (9%)	10 (7%)	1	0
All	All	916/926 (99%)	836 (91%)	56 (6%)	24 (3%)	4	1

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	SER
1	A	84	LYS
1	A	270	VAL
2	B	132	ASN
1	C	75	ASP
1	C	270	VAL
2	D	3	HIS
2	D	4	ASP
2	D	8	GLN
2	D	10	GLU
2	D	12	ILE
2	D	86	ILE
2	D	89	TYR
1	A	75	ASP
2	B	6	LYS
1	A	78	ASN
2	B	53	MET
2	B	7	LEU
2	B	34	LYS
1	A	79	THR
1	A	219	MET
2	D	11	ALA
2	D	91	VAL
2	D	88	ASN

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	261/261 (100%)	236 (90%)	25 (10%)	8	5
1	C	261/261 (100%)	246 (94%)	15 (6%)	18	17
2	B	135/137 (98%)	121 (90%)	14 (10%)	7	4
2	D	135/137 (98%)	121 (90%)	14 (10%)	7	4
All	All	792/796 (100%)	724 (91%)	68 (9%)	10	7

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ILE
1	A	35	GLN
1	A	59	PHE
1	A	65	ARG
1	A	76	SER
1	A	80	SER
1	A	116	THR
1	A	124	VAL
1	A	140	LEU
1	A	151	ARG
1	A	187	ILE
1	A	198	ILE
1	A	218	VAL
1	A	224	ILE
1	A	231	GLN
1	A	233	GLU
1	A	236	ASP
1	A	243	VAL
1	A	246	GLN
1	A	256	ASN
1	A	272	GLU
1	A	285	TYR
1	A	293	ILE
1	A	307	ASP
1	A	308	LEU
2	B	4	ASP
2	B	5	ASN
2	B	7	LEU
2	B	10	GLU
2	B	28	LYS
2	B	41	ARG
2	B	50	SER
2	B	52	GLU
2	B	58	LEU
2	B	85	ARG
2	B	96	ARG
2	B	133	ASP
2	B	144	GLU
2	B	153	ASN
1	C	37	GLU
1	C	59	PHE
1	C	74	SER
1	C	78	ASN

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Mol	Chain	Res	Type
1	C	81	LEU
1	C	140	LEU
1	C	145	ILE
1	C	167	ARG
1	C	198	ILE
1	C	232	LYS
1	C	250	ARG
1	C	270	VAL
1	C	272	GLU
1	C	279	LYS
1	C	285	TYR
2	D	4	ASP
2	D	7	LEU
2	D	8	GLN
2	D	10	GLU
2	D	14	ARG
2	D	60	LYS
2	D	68	GLU
2	D	80	GLN
2	D	85	ARG
2	D	86	ILE
2	D	94	LYS
2	D	101	GLU
2	D	151	LEU
2	D	153	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	21	ASN
1	A	156	HIS
1	A	196	GLN
1	A	212	HIS
1	A	256	ASN
1	A	291	ASN
1	A	297	GLN
1	A	305	ASN
2	B	8	GLN
2	B	40	GLN
2	B	73	GLN
2	B	84	ASN

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Mol	Chain	Res	Type
2	B	148	ASN
2	B	153	ASN
1	C	13	ASN
1	C	21	ASN
1	C	33	ASN
1	C	35	GLN
1	C	137	GLN
1	C	156	HIS
1	C	291	ASN
2	D	8	GLN
2	D	24	GLN
2	D	63	ASN
2	D	70	GLN
2	D	73	GLN
2	D	84	ASN
2	D	105	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 6 ligands modelled in this entry, 2 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	EOP	C	812	-	17,17,17	1.54	4 (23%)	20,24,24	0.56	0
5	CTP	D	902	-	29,30,30	1.60	5 (17%)	43,47,47	1.02	2 (4%)
5	CTP	B	901	-	29,30,30	1.61	7 (24%)	43,47,47	1.04	3 (6%)
3	EOP	A	811	-	17,17,17	1.54	4 (23%)	20,24,24	0.57	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	EOP	C	812	-	-	6/17/17/17	-
5	CTP	D	902	-	-	5/22/38/38	0/2/2/2
5	CTP	B	901	-	-	3/22/38/38	0/2/2/2
3	EOP	A	811	-	-	9/17/17/17	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	902	CTP	PG-O1G	3.52	1.61	1.50
5	B	901	CTP	PG-O1G	3.50	1.61	1.50
5	D	902	CTP	C4-N3	3.11	1.40	1.34
5	B	901	CTP	PB-O1B	3.07	1.61	1.50
5	D	902	CTP	PB-O1B	3.07	1.61	1.50
5	D	902	CTP	PA-O1A	3.03	1.61	1.50
5	B	901	CTP	PA-O1A	3.03	1.61	1.50
3	C	812	EOP	PAR-OAH	2.99	1.61	1.55
3	C	812	EOP	PAR-OAD	2.97	1.61	1.55
5	B	901	CTP	C4-N3	2.96	1.40	1.34
3	A	811	EOP	PAQ-OAE	2.96	1.61	1.55
3	C	812	EOP	PAQ-OAF	2.94	1.61	1.55
3	A	811	EOP	PAR-OAH	2.94	1.61	1.55
3	A	811	EOP	PAQ-OAF	2.93	1.61	1.55
3	A	811	EOP	PAR-OAD	2.92	1.61	1.55
3	C	812	EOP	PAQ-OAE	2.90	1.61	1.55
5	B	901	CTP	C2-N3	2.17	1.40	1.36
5	B	901	CTP	PA-O3A	2.06	1.61	1.59
5	D	902	CTP	C2-N3	2.05	1.40	1.36
5	B	901	CTP	PB-O3A	2.02	1.61	1.59

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	902	CTP	O3G-PG-O3B	2.68	113.64	104.64
5	B	901	CTP	O3G-PG-O3B	2.56	113.20	104.64
5	D	902	CTP	O2B-PB-O3A	2.40	113.77	107.27
5	B	901	CTP	O2B-PB-O3A	2.30	113.48	107.27
5	B	901	CTP	O4'-C1'-C2'	-2.01	102.31	106.62

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	811	EOP	CAK-CAO-NAM-CAI
3	A	811	EOP	CAO-CAK-PAQ-OAE
3	C	812	EOP	CAP-CAL-PAR-OAD
3	C	812	EOP	CAK-CAO-NAM-CAI
5	B	901	CTP	C5'-O5'-PA-O1A
5	B	901	CTP	C5'-O5'-PA-O3A
5	D	902	CTP	C5'-O5'-PA-O1A
5	D	902	CTP	C5'-O5'-PA-O2A
5	D	902	CTP	C5'-O5'-PA-O3A
5	D	902	CTP	PB-O3A-PA-O5'
3	A	811	EOP	OAA-CAO-NAM-CAI
3	C	812	EOP	OAA-CAO-NAM-CAI
3	A	811	EOP	NAM-CAI-CAJ-NAN
3	A	811	EOP	CAP-CAL-PAR-OAD
3	A	811	EOP	CAO-CAK-PAQ-OAF
3	C	812	EOP	CAP-CAL-PAR-OAH
3	C	812	EOP	NAM-CAI-CAJ-NAN
5	D	902	CTP	O4'-C4'-C5'-O5'
5	B	901	CTP	C5'-O5'-PA-O2A
3	A	811	EOP	CAP-CAL-PAR-OAH
3	A	811	EOP	CAP-CAL-PAR-OAG
3	A	811	EOP	CAO-CAK-PAQ-OAC
3	C	812	EOP	CAP-CAL-PAR-OAG

There are no ring outliers.

4 monomers are involved in 26 short contacts:

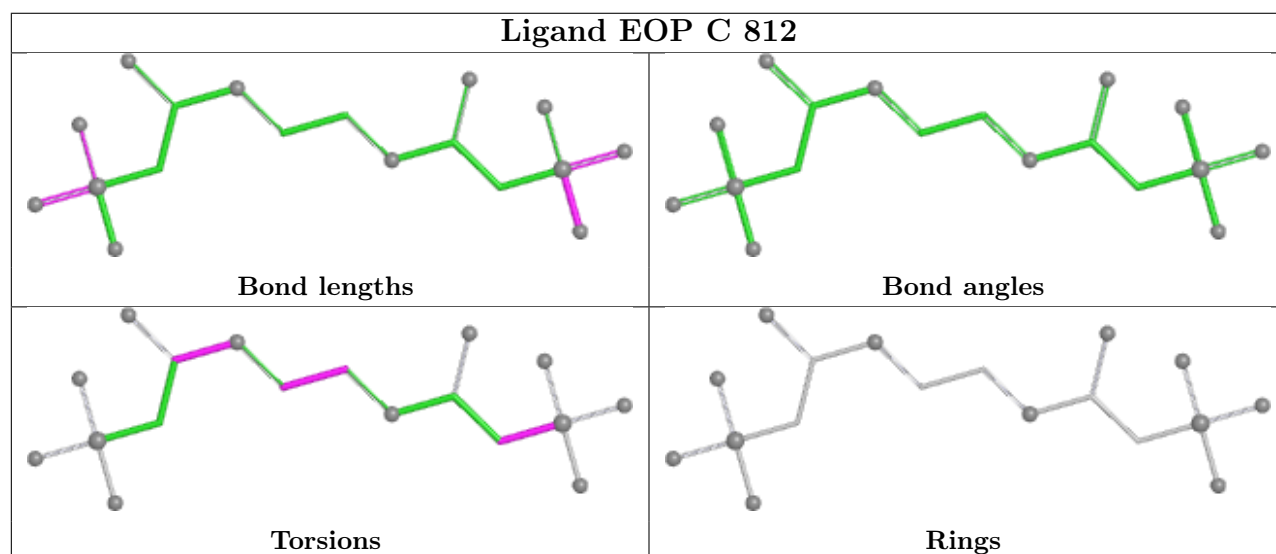
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	812	EOP	8	0
5	D	902	CTP	9	0

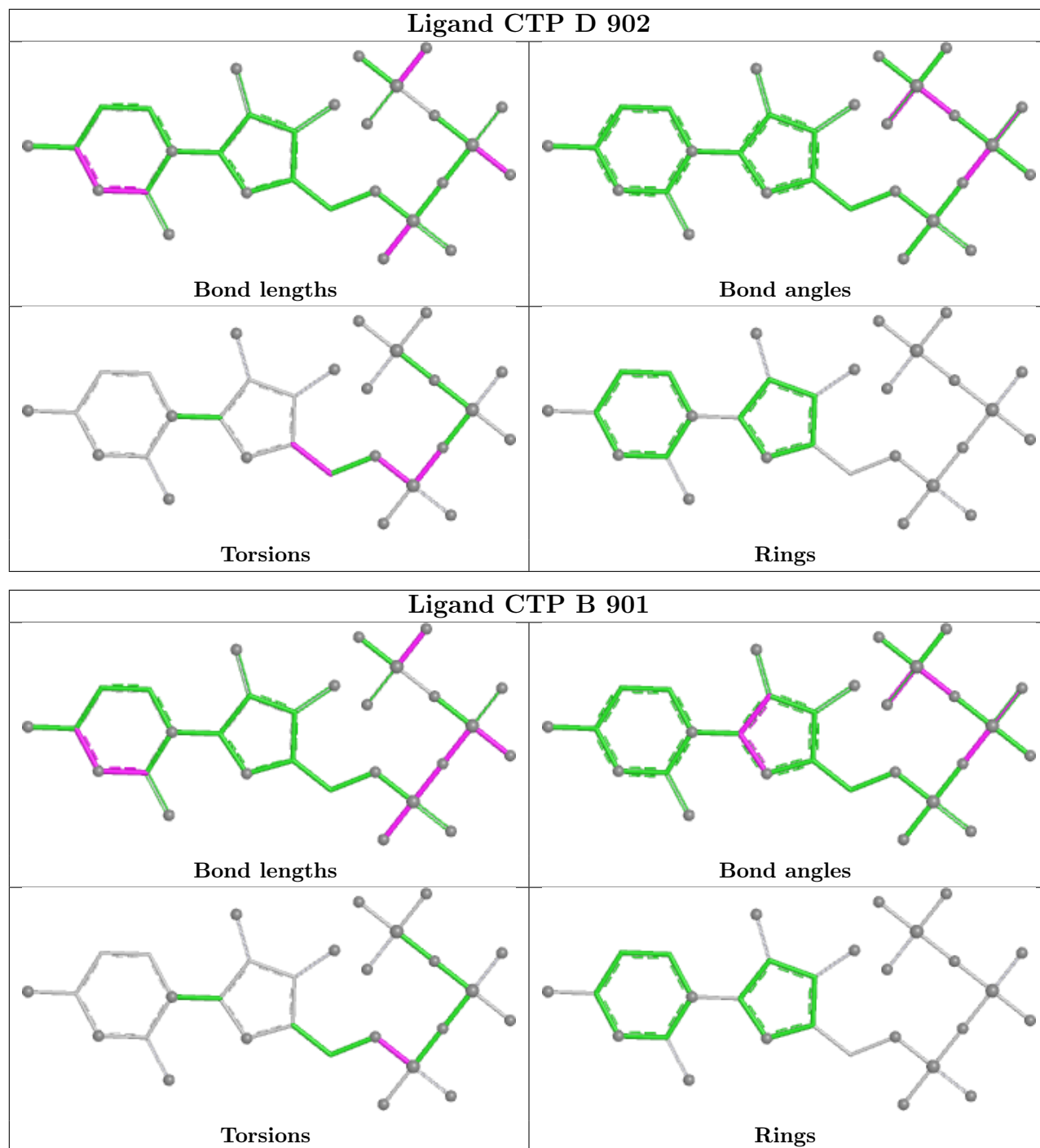
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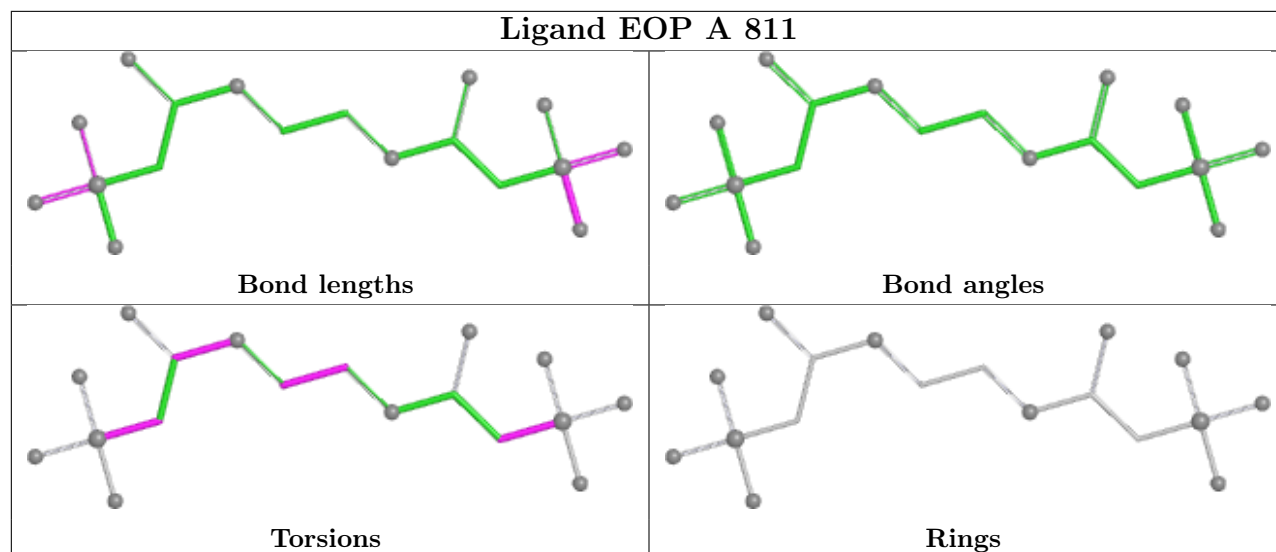
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	901	CTP	4	0
3	A	811	EOP	5	0

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.







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	310/310 (100%)	0.06	18 (5%)	29 30	19, 31, 64, 108	0
1	C	310/310 (100%)	-0.17	14 (4%)	38 40	17, 27, 63, 99	0
2	B	152/153 (99%)	1.00	25 (16%)	4 4	23, 50, 95, 125	0
2	D	152/153 (99%)	0.79	23 (15%)	5 5	24, 46, 97, 125	0
All	All	924/926 (99%)	0.26	80 (8%)	16 16	17, 34, 77, 125	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	86	ILE	9.5
1	A	81	LEU	8.2
1	A	82	GLY	6.5
1	A	80	SER	5.9
1	A	247	PHE	5.9
1	C	79	THR	5.7
1	C	81	LEU	5.5
2	B	7	LEU	5.4
1	C	80	SER	5.3
2	D	9	VAL	5.3
2	B	2	THR	4.8
2	B	12	ILE	4.6
1	A	77	ALA	4.5
2	D	12	ILE	4.5
2	B	9	VAL	4.5
2	D	11	ALA	4.4
1	C	78	ASN	4.3
1	A	308	LEU	4.3
2	D	2	THR	4.2
2	D	7	LEU	3.9
1	A	309	VAL	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	82	GLY	3.9
1	A	310	LEU	3.9
1	A	79	THR	3.9
2	B	11	ALA	3.9
1	C	247	PHE	3.7
1	A	85	GLY	3.7
2	B	3	HIS	3.7
2	B	82	THR	3.5
2	D	51	GLY	3.3
1	C	309	VAL	3.3
2	B	53	MET	3.3
1	C	235	LEU	3.3
2	B	8	GLN	3.3
2	D	92	VAL	3.2
2	B	92	VAL	3.2
2	D	53	MET	3.2
1	A	270	VAL	3.2
2	B	131	ALA	3.1
1	C	308	LEU	3.1
2	B	61	ILE	3.1
1	A	75	ASP	3.1
2	B	51	GLY	2.9
1	C	77	ALA	2.9
1	A	78	ASN	2.9
2	D	6	LYS	2.9
1	C	243	VAL	2.8
1	A	293	ILE	2.8
2	B	10	GLU	2.8
1	C	231	GLN	2.7
2	D	8	GLN	2.6
2	D	5	ASN	2.6
1	C	310	LEU	2.6
2	D	90	GLU	2.6
2	B	71	VAL	2.5
1	A	74	SER	2.5
1	A	76	SER	2.5
2	B	4	ASP	2.5
2	B	99	LEU	2.5
2	D	91	VAL	2.4
2	B	13	LYS	2.4
2	B	38	THR	2.4
2	D	48	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	68	GLU	2.4
1	C	234	ARG	2.4
1	A	54	ARG	2.3
1	A	246	GLN	2.3
2	B	5	ASN	2.3
2	D	4	ASP	2.2
2	D	82	THR	2.2
2	D	54	GLY	2.2
2	B	84	ASN	2.1
2	B	97	PRO	2.1
2	B	54	GLY	2.1
2	D	55	ARG	2.1
2	D	101	GLU	2.0
2	D	46	LEU	2.0
2	B	41	ARG	2.0
2	D	14	ARG	2.0
2	D	153	ASN	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 ⓘ

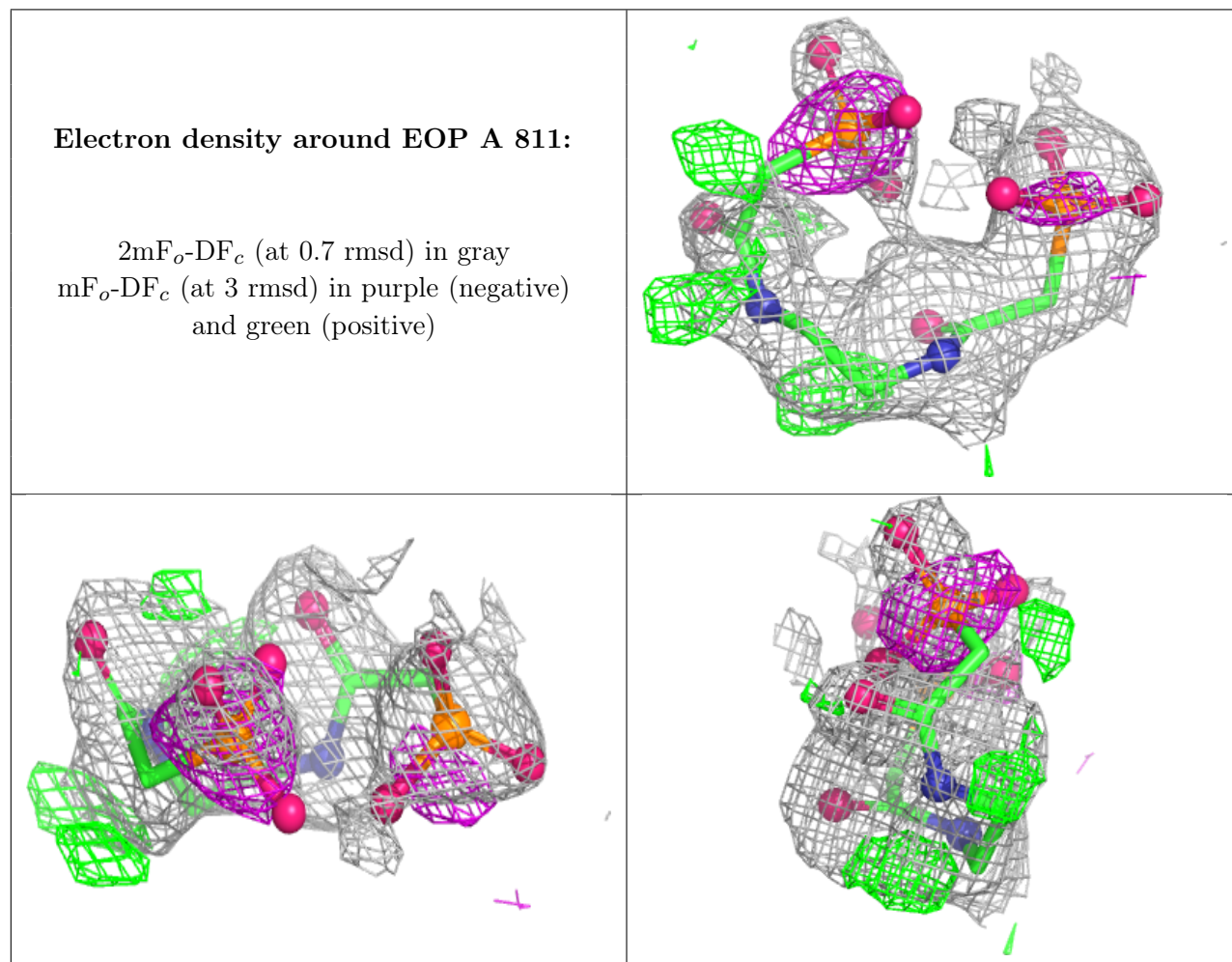
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

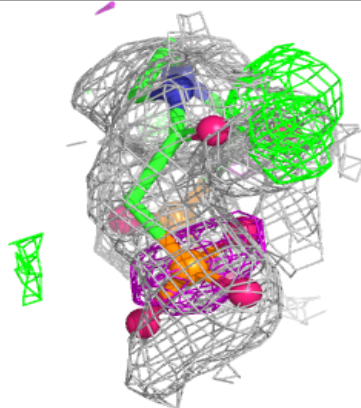
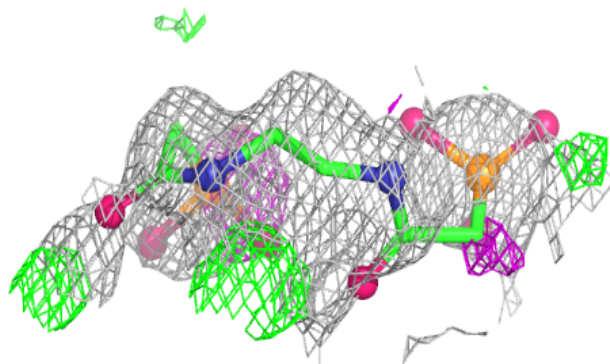
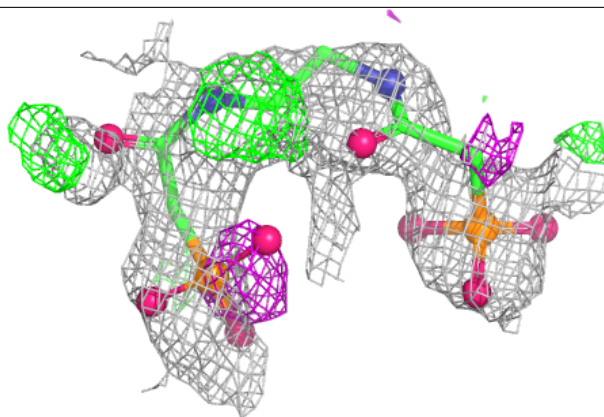
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	EOP	A	811	18/18	0.59	0.19	45,59,67,70	18
3	EOP	C	812	18/18	0.63	0.19	51,58,66,67	18
5	CTP	B	901	29/29	0.80	0.13	39,65,87,88	0
5	CTP	D	902	29/29	0.86	0.12	43,58,83,85	0
4	ZN	B	154	1/1	0.99	0.03	24,24,24,24	0
4	ZN	D	154	1/1	1.00	0.03	29,29,29,29	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.

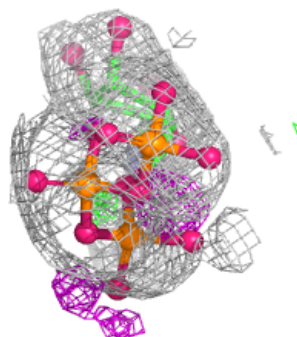
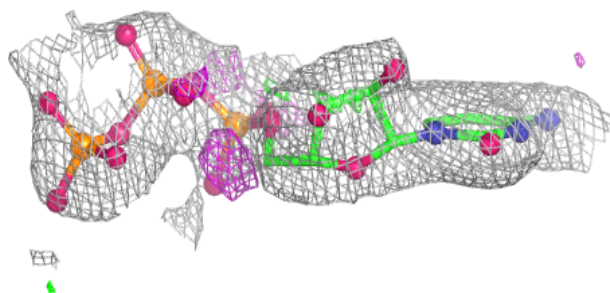
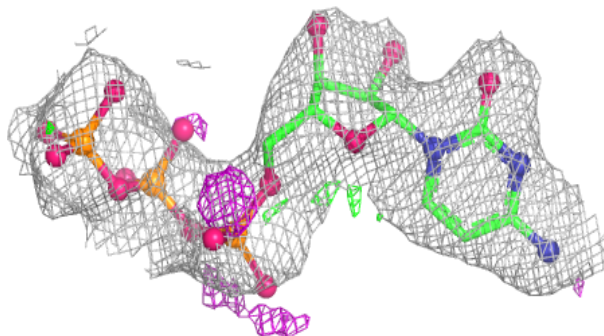


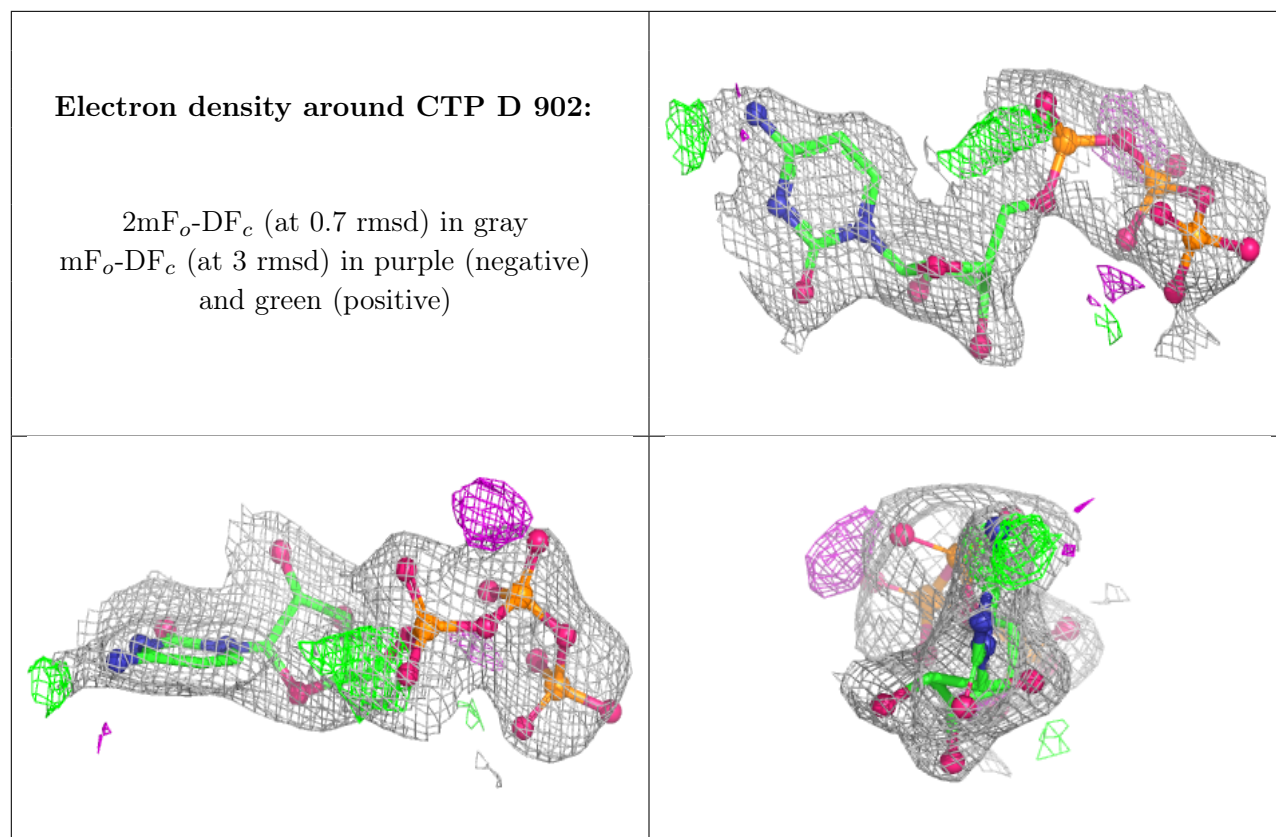
Electron density around EOP C 812:

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

**Electron density around CTP B 901:**

$2mF_o-DF_c$ (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.