



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:42 PM UTC

PDB ID : 5EXD / pdb_00005exd
Title : Crystal structure of oxalate oxidoreductase from *Moorella thermoacetica* bound with carboxy-di-oxido-methyl-TPP (COOM-TPP) intermediate
Authors : Gibson, M.I.; Chen, P.Y.-T.; Drennan, C.L.
Deposited on : 2015-11-23
Resolution : 2.50 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

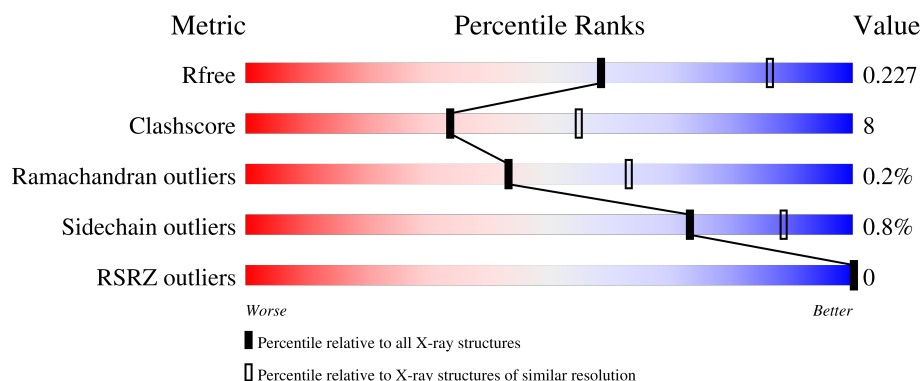
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	395	
1	D	395	
1	G	395	
1	J	395	
2	B	315	

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Mol	Chain	Length	Quality of chain
2	E	315	 78%16%6%
2	H	315	 76%17%8%
2	K	315	 74%15%11%
3	C	314	 82%17%..
3	F	314	 81%18%.
3	I	314	 79%18%..
3	L	314	 75%24%..

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
4	SF4	C	401	-	-	X	-
4	SF4	F	401	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 30592 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 Oxalate oxidoreductase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3060	1948	522	576	14			
1	D	394	Total	C	N	O	S	0	0	0
			3055	1944	523	574	14			
1	G	394	Total	C	N	O	S	0	0	0
			3058	1945	523	576	14			
1	J	394	Total	C	N	O	S	0	0	0
			3058	1947	521	576	14			

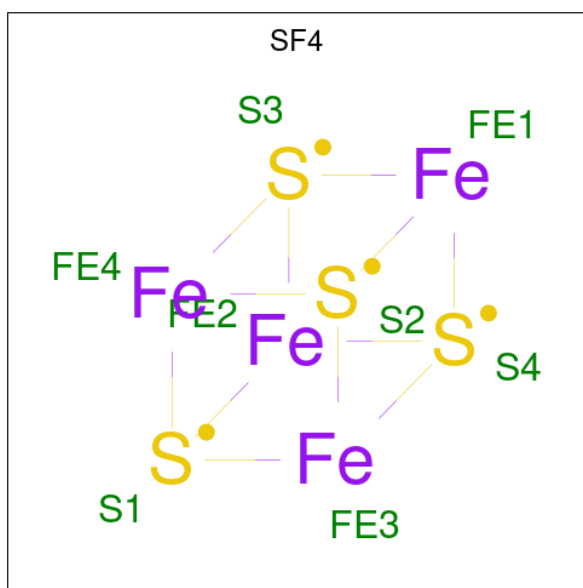
- Molecule 2 is a protein called Oxalate oxidoreductase subunit delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	286	Total	C	N	O	S	0	0	0
			2162	1369	361	417	15			
2	E	295	Total	C	N	O	S	0	0	0
			2180	1369	368	429	14			
2	H	291	Total	C	N	O	S	0	0	0
			2200	1390	371	424	15			
2	K	280	Total	C	N	O	S	0	0	0
			2091	1317	352	407	15			

- Molecule 3 is a protein called Oxalate oxidoreductase subunit beta.

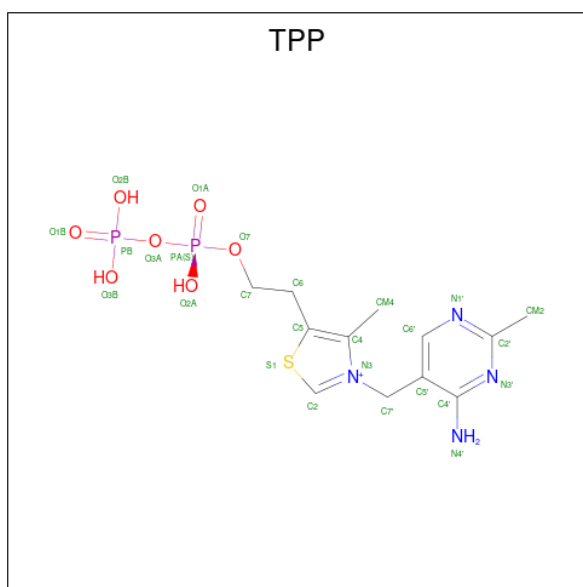
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	312	Total	C	N	O	S	0	0	0
			2393	1535	409	432	17			
3	F	310	Total	C	N	O	S	0	0	0
			2371	1522	405	427	17			
3	I	309	Total	C	N	O	S	0	0	0
			2362	1516	402	428	16			
3	L	310	Total	C	N	O	S	0	0	0
			2375	1524	405	429	17			

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	Fe	S	0	0
			8	4	4		
4	B	1	Total	Fe	S	0	0
			8	4	4		
4	C	1	Total	Fe	S	0	0
			8	4	4		
4	E	1	Total	Fe	S	0	0
			8	4	4		
4	E	1	Total	Fe	S	0	0
			8	4	4		
4	F	1	Total	Fe	S	0	0
			8	4	4		
4	H	1	Total	Fe	S	0	0
			8	4	4		
4	H	1	Total	Fe	S	0	0
			8	4	4		
4	I	1	Total	Fe	S	0	0
			8	4	4		
4	K	1	Total	Fe	S	0	0
			8	4	4		
4	K	1	Total	Fe	S	0	0
			8	4	4		
4	L	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 5 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: $\text{C}_{12}\text{H}_{19}\text{N}_4\text{O}_7\text{P}_2\text{S}$).

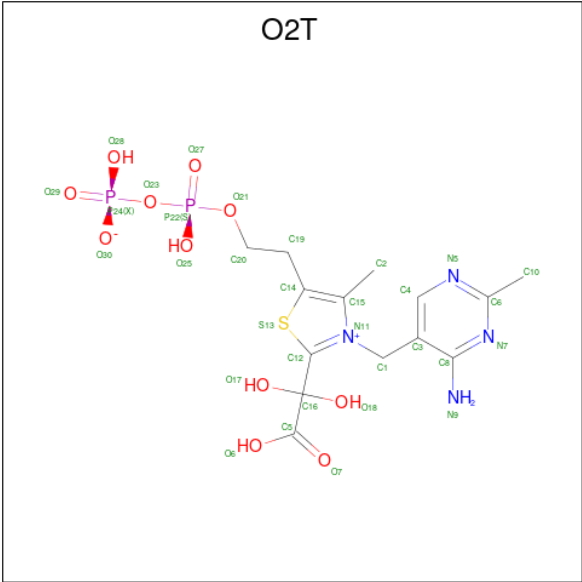


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	C	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		
5	I	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		
5	L	1	Total	C	N	O	P	S	0	0
			26	12	4	7	2	1		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Mg	0	0
			1	1		
6	F	1	Total	Mg	0	0
			1	1		
6	I	1	Total	Mg	0	0
			1	1		
6	L	1	Total	Mg	0	0
			1	1		

- Molecule 7 is [2-[3-[(4-azanyl-2-methyl-pyrimidin-5-yl)methyl]-4-methyl-2-[1,1,2-tris(oxidanyl)-2-oxidanylidene-ethyl]-1,3-thiazol-3-ium-5-yl]ethoxy-oxidanyl-phosphoryl] hydrogen phosphate (CCD ID: O2T) (formula: C₁₄H₂₀N₄O₁₁P₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
7	F	1	Total	C	N	O	P	S	0	1
			32	14	4	11	2	1		

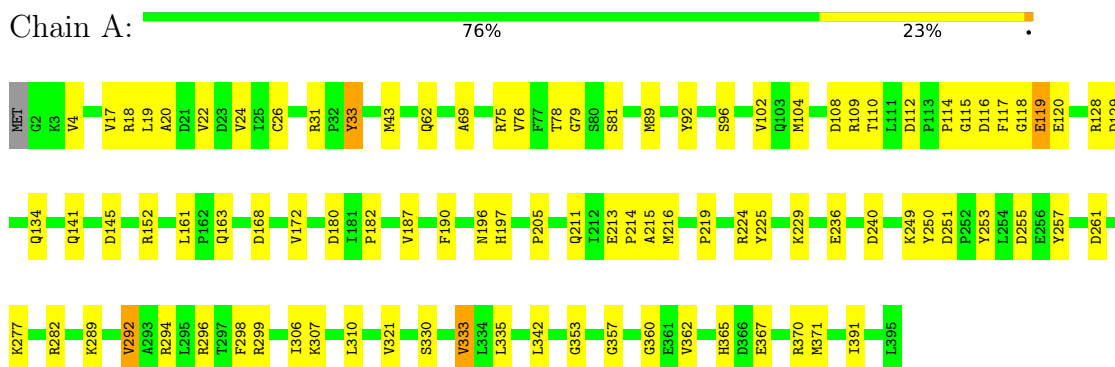
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	O	0	0
			3	3		
8	C	1	Total	O	0	0
			1	1		
8	D	1	Total	O	0	0
			1	1		
8	F	4	Total	O	0	0
			4	4		
8	G	2	Total	O	0	0
			2	2		
8	H	1	Total	O	0	0
			1	1		
8	I	4	Total	O	0	0
			4	4		
8	L	1	Total	O	0	0
			1	1		

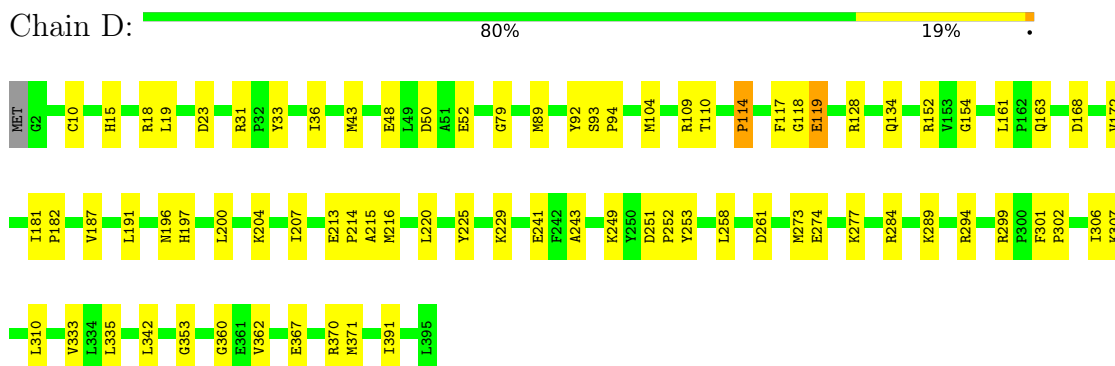
3 Residue-property plots [i](#)

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.

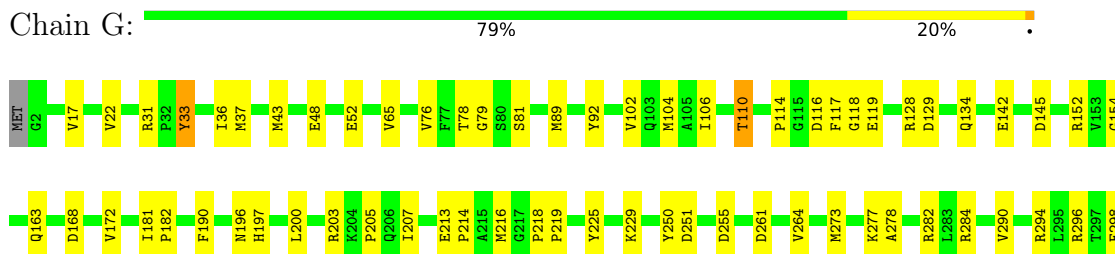
- Molecule 1: Oxalate oxidoreductase subunit alpha



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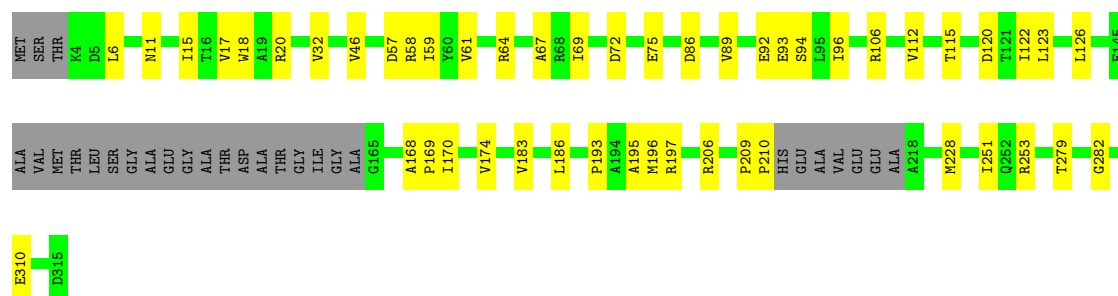
- Molecule 1: Oxalate oxidoreductase subunit alpha

Chain J: 79% 20%



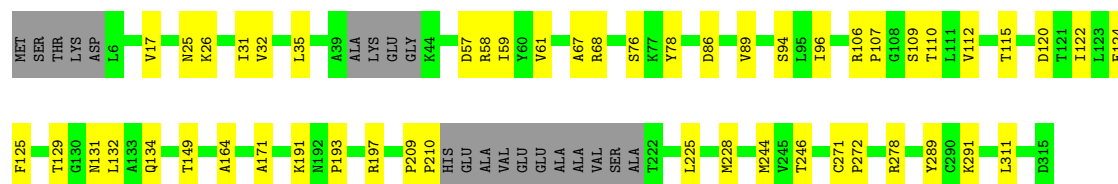
- Molecule 2: Oxalate oxidoreductase subunit delta

Chain B: 75% 16% 9%



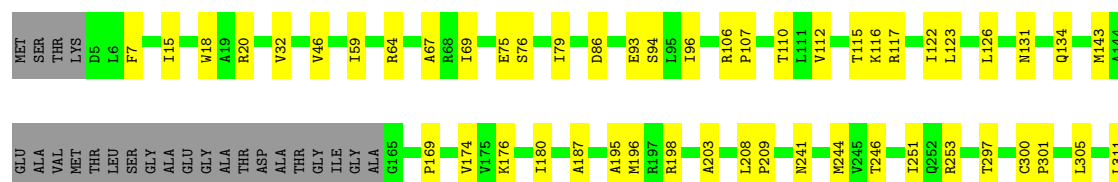
- Molecule 2: Oxalate oxidoreductase subunit delta

Chain E: 78% 16% 6%



- Molecule 2: Oxalate oxidoreductase subunit delta

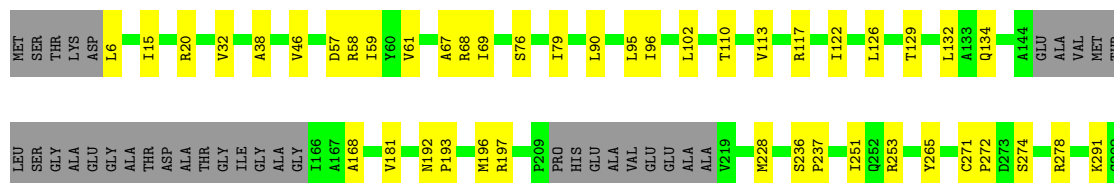
Chain H: 76% 17% 8%





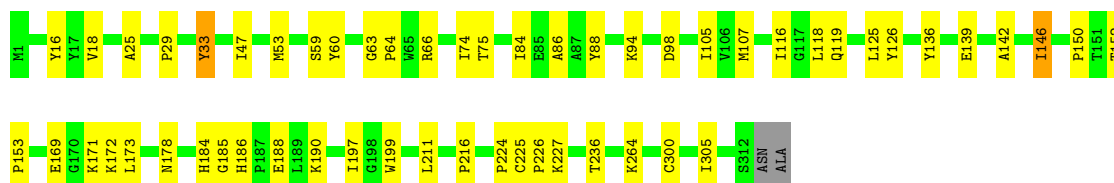
- Molecule 2: Oxalate oxidoreductase subunit delta

Chain K: 74% 15% 11%



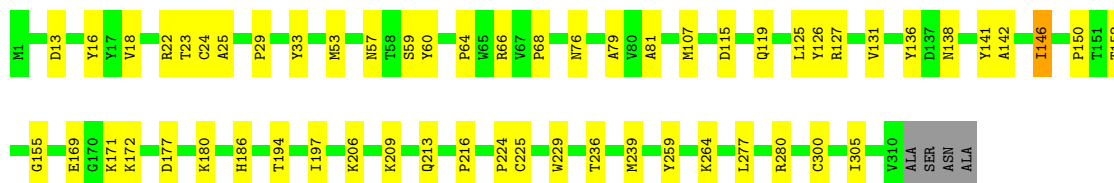
- Molecule 3: Oxalate oxidoreductase subunit beta

Chain C: 82% 17% ..



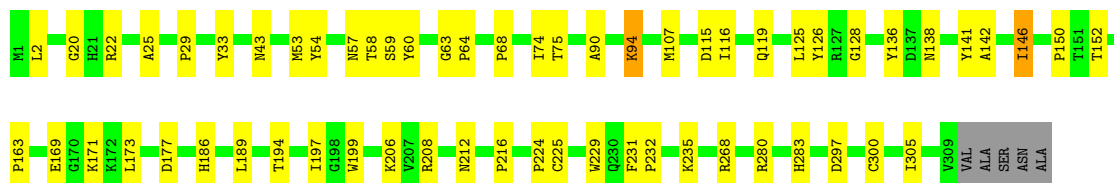
- Molecule 3: Oxalate oxidoreductase subunit beta

Chain F: 81% 18% .



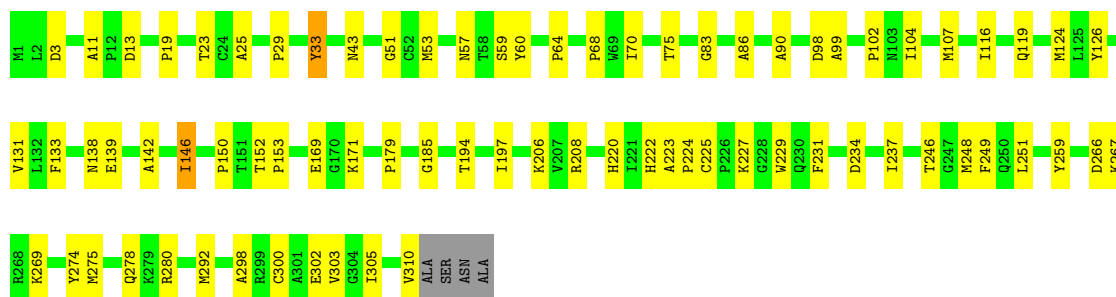
- Molecule 3: Oxalate oxidoreductase subunit beta

Chain I: 79% 18% ..



- Molecule 3: Oxalate oxidoreductase subunit beta

Chain L: 75% 24% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	138.43Å 138.43Å 217.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.94 – 2.50 48.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.2 (48.94-2.50) 93.5 (48.94-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.196 , 0.229 0.196 , 0.227	Depositor DCC
R_{free} test set	6540 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.947	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 7.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.328 for h,-k,-l	Xtriage
Reported twinning fraction	0.640 for h,-k,-l	Depositor
Outliers	0 of 131805 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	30592	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.71% 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: MG, O2T, SF4, TPP

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	0.31	0/3132	0.74	0/4248
1	D	0.31	0/3127	0.75	4/4243 (0.1%)
1	G	0.32	0/3130	0.74	0/4247
1	J	0.31	0/3130	0.73	2/4246 (0.0%)
2	B	0.31	0/2201	0.73	0/3003
2	E	0.31	0/2219	0.76	0/3036
2	H	0.31	0/2241	0.77	3/3059 (0.1%)
2	K	0.29	0/2129	0.75	0/2911
3	C	0.33	0/2458	0.76	2/3339 (0.1%)
3	F	0.32	0/2436	0.76	0/3312
3	I	0.32	0/2427	0.74	2/3300 (0.1%)
3	L	0.31	0/2440	0.77	2/3317 (0.1%)
All	All	0.31	0/31070	0.75	15/42261 (0.0%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	208	LEU	CA-C-N	6.20	124.19	119.66
2	H	208	LEU	C-N-CA	6.20	124.19	119.66
1	D	204	LYS	CA-C-N	5.76	125.52	119.76
1	D	204	LYS	C-N-CA	5.76	125.52	119.76
3	C	199	TRP	CA-C-N	5.55	124.90	118.85
3	C	199	TRP	C-N-CA	5.55	124.90	118.85
1	D	191	LEU	CA-C-N	5.41	125.95	120.38
1	D	191	LEU	C-N-CA	5.41	125.95	120.38
3	L	223	ALA	CA-C-N	5.21	125.21	119.90
3	L	223	ALA	C-N-CA	5.21	125.21	119.90
1	J	204	LYS	CA-C-N	5.08	125.00	119.76
1	J	204	LYS	C-N-CA	5.08	125.00	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	209	PRO	O-C-N	5.04	123.63	121.31
3	I	199	TRP	CA-C-N	5.03	124.63	119.05
3	I	199	TRP	C-N-CA	5.03	124.63	119.05

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	3060	0	2992	62	0
1	D	3055	0	2979	50	0
1	G	3058	0	2981	64	0
1	J	3058	0	2985	56	0
2	B	2162	0	2149	30	0
2	E	2180	0	2102	33	0
2	H	2200	0	2182	31	0
2	K	2091	0	2019	35	0
3	C	2393	0	2394	40	0
3	F	2371	0	2360	42	0
3	I	2362	0	2346	44	0
3	L	2375	0	2364	50	0
4	B	16	0	0	0	0
4	C	8	0	0	2	0
4	E	16	0	0	0	0
4	F	8	0	0	2	0
4	H	16	0	0	0	0
4	I	8	0	0	1	0
4	K	16	0	0	0	0
4	L	8	0	0	1	0
5	C	26	0	16	1	0
5	I	26	0	16	1	0
5	L	26	0	16	1	0
6	C	1	0	0	0	0
6	F	1	0	0	0	0
6	I	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	1	0	0	0	0
7	F	32	0	2	2	0
8	A	3	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
8	F	4	0	0	1	0
8	G	2	0	0	0	0
8	H	1	0	0	0	0
8	I	4	0	0	0	0
8	L	1	0	0	0	0
All	All	30592	0	29903	465	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 8.

All (465) 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:G:216:MET:HE3	2:K:228:MET:HE2	1.59	0.84
2:E:164:ALA:HA	2:E:191:LYS:HB2	1.67	0.76
1:G:261:ASP:OD1	1:G:284:ARG:NH2	2.18	0.76
2:K:117:ARG:NH1	2:K:312:ASP:OD1	2.20	0.74
1:G:284:ARG:NH1	1:G:290:VAL:O	2.21	0.74
1:D:118:GLY:HA3	1:D:360:GLY:HA3	1.68	0.73
3:I:208:ARG:O	3:I:212:ASN:ND2	2.21	0.73
1:A:216:MET:HE3	2:E:228:MET:HE2	1.69	0.73
2:B:228:MET:HE2	1:D:216:MET:HE3	1.70	0.73
2:B:183:VAL:HG13	2:B:196:MET:HE1	1.73	0.71
3:L:152:THR:O	3:L:171:LYS:NZ	2.24	0.71
3:L:194:THR:O	3:L:206:LYS:NZ	2.23	0.71
1:A:129:ASP:OD2	1:D:128:ARG:NH1	2.24	0.70
3:C:94:LYS:NZ	1:D:50:ASP:O	2.18	0.69
3:C:119:GLN:HB2	3:F:119:GLN:HB2	1.76	0.68
3:F:194:THR:O	3:F:206:LYS:NZ	2.23	0.68
2:H:15:ILE:HB	2:H:67:ALA:HB3	1.75	0.68
2:B:15:ILE:HB	2:B:67:ALA:HB3	1.74	0.67
3:F:300:CYS:HB3	3:F:305:ILE:HB	1.75	0.67
1:J:118:GLY:HA3	1:J:360:GLY:HA3	1.76	0.67
1:J:19:LEU:HB3	1:J:187:VAL:HG21	1.76	0.67
3:I:119:GLN:HB2	3:L:119:GLN:HB2	1.77	0.66
3:I:300:CYS:HB3	3:I:305:ILE:HB	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:29:PRO:HB2	3:I:53:MET:HE3	1.79	0.65
2:H:69:ILE:HD12	2:H:180:ILE:HD11	1.79	0.65
1:A:277:LYS:HG2	1:A:292:VAL:HG21	1.78	0.64
3:C:94:LYS:HZ1	1:D:23:ASP:HA	1.60	0.64
2:E:25:ASN:ND2	2:E:57:ASP:OD2	2.31	0.64
1:A:96:SER:OG	1:A:224:ARG:NH1	2.31	0.64
1:A:255:ASP:HB3	1:A:294:ARG:HB3	1.79	0.64
1:G:118:GLY:HA3	1:G:360:GLY:HA3	1.78	0.64
1:A:110:THR:HB	1:A:114:PRO:HD2	1.80	0.63
2:K:96:ILE:HD13	2:K:122:ILE:HG23	1.79	0.63
1:A:240:ASP:OD1	1:A:249:LYS:NZ	2.32	0.62
1:A:257:TYR:HB3	1:A:292:VAL:HG23	1.80	0.62
2:K:68:ARG:NH1	2:K:76:SER:O	2.32	0.62
3:L:300:CYS:HB3	3:L:305:ILE:HB	1.82	0.62
1:A:89:MET:HE3	1:D:89:MET:HG2	1.81	0.61
2:B:112:VAL:HG21	2:B:174:VAL:HG22	1.82	0.61
2:H:20:ARG:HG3	2:H:251:ILE:HD11	1.82	0.61
2:B:168:ALA:HB1	2:B:196:MET:HB2	1.81	0.61
1:G:145:ASP:OD2	1:G:250:TYR:OH	2.18	0.61
2:E:59:ILE:HB	3:F:25:ALA:H	1.66	0.61
1:A:20:ALA:O	1:A:75:ARG:NH1	2.30	0.60
3:C:185:GLY:HA2	3:F:177:ASP:HB3	1.83	0.60
2:E:291:LYS:NZ	3:F:23:THR:O	2.31	0.60
1:G:114:PRO:HB3	1:J:212:ILE:HD11	1.81	0.60
3:C:29:PRO:HB2	3:C:53:MET:HE3	1.83	0.60
2:E:86:ASP:OD2	2:E:106:ARG:NH2	2.29	0.60
1:D:15:HIS:HD2	1:D:18:ARG:HH11	1.49	0.59
3:C:300:CYS:HB3	3:C:305:ILE:HB	1.85	0.59
3:F:142:ALA:HA	3:F:146:ILE:HD13	1.84	0.59
1:A:117:PHE:CE2	1:D:214:PRO:HB3	2.38	0.59
1:A:282:ARG:NH2	2:B:6:LEU:O	2.36	0.59
3:I:125:LEU:HD22	3:I:189:LEU:HD13	1.85	0.59
2:H:112:VAL:HG21	2:H:174:VAL:HG22	1.85	0.58
1:A:168:ASP:HB2	1:A:172:VAL:HG23	1.86	0.58
1:D:367:GLU:HG2	1:D:370:ARG:HH12	1.68	0.58
3:C:126:TYR:CZ	3:F:150:PRO:HA	2.38	0.58
3:F:29:PRO:HB2	3:F:53:MET:HE3	1.86	0.58
1:G:31:ARG:NH2	1:G:116:ASP:OD2	2.36	0.58
1:D:273:MET:HG3	1:D:294:ARG:HG3	1.85	0.58
1:A:128:ARG:HG2	1:A:298:PHE:HE2	1.69	0.57
2:E:32:VAL:HG13	2:E:67:ALA:HB2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:90:ALA:O	3:I:94:LYS:HD3	2.03	0.57
1:J:110:THR:HB	1:J:114:PRO:HD2	1.86	0.57
1:J:109:ARG:HD3	1:J:110:THR:O	2.03	0.57
1:G:168:ASP:HB2	1:G:172:VAL:HG23	1.86	0.57
2:K:253:ARG:NH2	2:K:310:GLU:OE1	2.37	0.57
3:I:142:ALA:HA	3:I:146:ILE:HD13	1.87	0.57
3:F:24:CYS:HB2	4:F:401:SF4:S3	2.44	0.57
2:H:32:VAL:HG13	2:H:67:ALA:HB2	1.86	0.57
2:H:297:THR:HG22	2:H:305:LEU:HB3	1.86	0.57
2:B:20:ARG:HG3	2:B:251:ILE:HD11	1.87	0.56
1:D:110:THR:HB	1:D:114:PRO:HD2	1.86	0.56
1:D:258:LEU:HD13	1:D:284:ARG:HG3	1.86	0.56
1:J:282:ARG:NH1	2:K:6:LEU:O	2.36	0.56
1:A:145:ASP:OD2	1:A:250:TYR:OH	2.20	0.56
2:K:278:ARG:NH1	3:L:11:ALA:O	2.38	0.56
1:A:128:ARG:HD3	1:A:333:VAL:HG21	1.88	0.56
2:B:32:VAL:HG13	2:B:67:ALA:HB2	1.86	0.56
1:A:161:LEU:O	1:A:163:GLN:NE2	2.34	0.56
2:H:123:LEU:HD23	2:H:126:LEU:HD12	1.87	0.56
2:K:110:THR:HG23	2:K:134:GLN:HG3	1.88	0.55
2:H:169:PRO:HG3	2:H:195:ALA:HB1	1.88	0.55
1:A:31:ARG:NH1	1:A:116:ASP:O	2.39	0.55
2:E:193:PRO:O	2:E:197:ARG:HG2	2.07	0.55
1:G:128:ARG:HD3	1:G:333:VAL:HG21	1.88	0.55
3:I:58:THR:HG21	1:J:211:GLN:H	1.71	0.55
3:L:29:PRO:HB2	3:L:53:MET:HE3	1.88	0.55
3:F:152:THR:HB	3:F:171:LYS:HE3	1.89	0.55
3:L:246:THR:HG22	3:L:269:LYS:HB2	1.89	0.55
2:H:46:VAL:HG12	2:H:69:ILE:HG12	1.87	0.55
1:G:330:SER:OG	1:J:229:LYS:NZ	2.37	0.55
3:C:171:LYS:NZ	3:C:173:LEU:O	2.39	0.54
3:L:227:LYS:HD3	4:L:401:SF4:S2	2.48	0.54
2:H:59:ILE:HB	3:I:25:ALA:H	1.73	0.54
3:L:142:ALA:HA	3:L:146:ILE:HD13	1.88	0.54
1:A:196:ASN:OD1	1:A:197:HIS:N	2.40	0.54
1:A:330:SER:OG	1:D:229:LYS:NZ	2.41	0.54
2:B:123:LEU:HD23	2:B:126:LEU:HD12	1.90	0.54
2:K:96:ILE:HD11	2:K:113:VAL:HG11	1.90	0.54
1:G:89:MET:HA	1:G:92:TYR:CE2	2.43	0.53
1:G:205:PRO:HB2	2:K:236:SER:HB2	1.89	0.53
2:K:15:ILE:HB	2:K:67:ALA:HB3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:32:VAL:HG13	2:K:67:ALA:HB2	1.90	0.53
3:C:224:PRO:HB2	3:C:236:THR:HG23	1.91	0.53
3:C:16:TYR:HA	3:C:66:ARG:HD3	1.90	0.53
1:D:367:GLU:HG2	1:D:370:ARG:NH1	2.22	0.53
2:E:129:THR:HB	2:E:132:LEU:HB3	1.91	0.53
1:J:258:LEU:HD13	1:J:284:ARG:HG3	1.88	0.53
1:J:145:ASP:OD2	1:J:250:TYR:OH	2.22	0.53
3:C:142:ALA:HA	3:C:146:ILE:HD13	1.91	0.53
1:J:20:ALA:O	1:J:75:ARG:NH1	2.38	0.53
1:J:134:GLN:HB3	1:J:299:ARG:HB2	1.91	0.53
1:A:205:PRO:HA	3:F:68:PRO:HG3	1.91	0.52
2:K:168:ALA:HB1	2:K:196:MET:HB2	1.90	0.52
3:L:274:TYR:O	3:L:278:GLN:NE2	2.41	0.52
1:A:79:GLY:HA2	1:A:104:MET:HG3	1.92	0.52
3:L:75:THR:HG22	3:L:116:ILE:HG12	1.91	0.52
3:C:94:LYS:NZ	1:D:23:ASP:HA	2.24	0.52
2:H:143:MET:HG2	2:H:198:ARG:HE	1.74	0.52
2:K:46:VAL:HG12	2:K:69:ILE:HG12	1.91	0.52
3:L:33:TYR:CE1	3:L:107:MET:HG3	2.44	0.52
3:C:188:GLU:O	3:C:190:LYS:NZ	2.42	0.52
2:K:57:ASP:OD1	2:K:57:ASP:N	2.43	0.52
3:I:150:PRO:HA	3:L:126:TYR:CZ	2.45	0.52
3:I:171:LYS:NZ	3:I:173:LEU:O	2.42	0.51
1:D:10:CYS:HA	1:D:36:ILE:HG12	1.91	0.51
2:B:228:MET:HE3	1:D:220:LEU:HB2	1.92	0.51
1:D:19:LEU:HB3	1:D:187:VAL:HG21	1.92	0.51
3:C:75:THR:HG22	3:C:116:ILE:HG12	1.93	0.51
3:L:138:ASN:HB2	3:L:225:CYS:HB2	1.93	0.51
1:G:79:GLY:HA2	1:G:104:MET:HG3	1.92	0.51
2:E:26:LYS:HD2	2:E:149:THR:HG22	1.92	0.51
1:G:255:ASP:OD2	1:G:277:LYS:NZ	2.34	0.51
2:H:76:SER:HB2	2:H:79:ILE:HD12	1.91	0.51
3:I:177:ASP:HB3	3:L:185:GLY:HA2	1.93	0.51
3:I:268:ARG:NH2	3:I:297:ASP:OD1	2.37	0.51
1:A:251:ASP:O	1:A:296:ARG:NH1	2.43	0.50
3:F:127:ARG:NH2	8:F:502:HOH:O	2.28	0.50
3:I:197:ILE:HD11	3:I:224:PRO:HD2	1.94	0.50
1:J:196:ASN:OD1	1:J:197:HIS:N	2.39	0.50
1:G:89:MET:HE3	1:J:89:MET:HG2	1.94	0.50
3:I:229:TRP:HA	3:I:280:ARG:HD3	1.93	0.50
2:B:92:GLU:OE1	2:B:94:SER:OG	2.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:104:MET:HE2	1:G:106:ILE:HG12	1.94	0.50
2:B:169:PRO:HG3	2:B:195:ALA:HB1	1.93	0.50
1:G:134:GLN:HB3	1:G:299:ARG:HB2	1.93	0.50
3:I:194:THR:O	3:I:206:LYS:NZ	2.37	0.50
1:A:214:PRO:HB3	1:D:117:PHE:CE2	2.47	0.50
2:H:253:ARG:NH2	2:H:315:ASP:OXT	2.35	0.50
2:E:68:ARG:NH1	2:E:76:SER:OG	2.45	0.50
2:K:59:ILE:HB	3:L:25:ALA:H	1.76	0.50
3:C:125:LEU:HD23	3:C:216:PRO:HB2	1.94	0.49
1:A:367:GLU:O	1:A:371:MET:HG3	2.12	0.49
2:E:57:ASP:OD1	2:E:57:ASP:N	2.45	0.49
1:A:134:GLN:HB3	1:A:299:ARG:HB2	1.94	0.49
1:J:255:ASP:HB3	1:J:294:ARG:HB3	1.93	0.49
3:L:197:ILE:HD11	3:L:224:PRO:HD2	1.94	0.49
1:G:225:TYR:O	1:G:229:LYS:HG2	2.13	0.49
1:J:371:MET:HE2	1:J:391:ILE:HD12	1.94	0.49
3:C:184:HIS:HB3	3:F:180:LYS:HB3	1.94	0.49
1:G:251:ASP:O	1:G:296:ARG:HD3	2.12	0.49
2:H:241:ASN:HB3	3:I:20:GLY:HA3	1.95	0.49
1:A:19:LEU:HB3	1:A:187:VAL:HG21	1.95	0.49
1:G:306:ILE:HG23	1:G:310:LEU:HD12	1.95	0.49
1:G:367:GLU:HG2	1:G:370:ARG:NH2	2.28	0.49
2:K:20:ARG:HG3	2:K:251:ILE:HD11	1.94	0.49
1:D:306:ILE:HG23	1:D:310:LEU:HD12	1.95	0.49
1:A:367:GLU:HG2	1:A:370:ARG:NH2	2.27	0.49
3:F:239:MET:HE1	3:F:277:LEU:HB3	1.94	0.49
1:G:255:ASP:OD2	1:G:294:ARG:NH2	2.46	0.49
1:J:76:VAL:HG21	1:J:102:VAL:HG22	1.95	0.49
1:G:196:ASN:OD1	1:G:197:HIS:N	2.43	0.48
3:C:118:LEU:HD13	3:C:178:ASN:HD21	1.77	0.48
1:G:214:PRO:HB3	1:J:117:PHE:CZ	2.48	0.48
1:A:24:VAL:HB	1:A:76:VAL:HG12	1.96	0.48
1:G:129:ASP:OD2	1:J:128:ARG:NH2	2.46	0.48
3:L:275:MET:SD	3:L:292:MET:HG2	2.54	0.48
1:A:19:LEU:HD13	1:A:182:PRO:HG2	1.95	0.48
3:F:33:TYR:CE1	3:F:107:MET:HG3	2.48	0.48
1:G:282:ARG:NH1	2:H:7:PHE:HA	2.28	0.48
1:A:118:GLY:HA3	1:A:360:GLY:HA3	1.95	0.48
1:A:214:PRO:HB3	1:D:117:PHE:CZ	2.49	0.48
1:D:152:ARG:HH22	1:D:241:GLU:HG3	1.78	0.48
3:F:115:ASP:OD2	3:F:141:TYR:OH	2.30	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:321:VAL:HG13	1:J:357:GLY:HA3	1.94	0.48
2:K:102:LEU:HD11	2:K:126:LEU:HD22	1.95	0.48
1:A:109:ARG:HA	1:A:119:GLU:HA	1.95	0.48
1:G:255:ASP:HB3	1:G:294:ARG:HB3	1.95	0.48
1:G:278:ALA:O	1:G:282:ARG:HG3	2.14	0.48
3:C:150:PRO:HA	3:F:126:TYR:CZ	2.49	0.47
1:J:367:GLU:O	1:J:371:MET:HG3	2.13	0.47
3:L:43:ASN:OD1	3:L:102:PRO:HA	2.14	0.47
3:L:124:MET:HE1	3:L:133:PHE:HB2	1.96	0.47
2:K:76:SER:HB2	2:K:79:ILE:HD12	1.96	0.47
3:L:59:SER:HA	3:L:60:TYR:HA	1.56	0.47
3:I:169:GLU:OE1	3:I:169:GLU:N	2.46	0.47
1:J:99:ARG:NH2	1:J:159:VAL:O	2.38	0.47
1:D:274:GLU:HB2	2:E:78:TYR:HE2	1.80	0.47
1:D:168:ASP:HB2	1:D:172:VAL:HG23	1.97	0.47
1:G:117:PHE:CE2	1:J:214:PRO:HB3	2.50	0.47
1:A:128:ARG:HG2	1:A:298:PHE:CE2	2.48	0.47
2:B:58:ARG:O	2:B:61:VAL:HG12	2.15	0.47
1:D:109:ARG:HA	1:D:119:GLU:HA	1.95	0.47
3:F:155:GLY:HA2	3:F:169:GLU:HB2	1.97	0.47
1:G:371:MET:HE2	1:G:391:ILE:HD12	1.97	0.47
2:H:187:ALA:HB2	2:H:196:MET:HE1	1.97	0.47
2:H:244:MET:HE2	3:I:22:ARG:HG3	1.97	0.47
1:J:335:LEU:HD22	1:J:353:GLY:HA3	1.96	0.47
1:A:236:GLU:OE2	1:A:253:TYR:OH	2.23	0.47
3:I:74:ILE:HD12	5:I:402:TPP:H72	1.96	0.47
3:L:98:ASP:OD1	3:L:98:ASP:N	2.48	0.47
1:A:216:MET:C	1:A:219:PRO:HD2	2.40	0.47
2:B:115:THR:HG21	2:B:122:ILE:HD11	1.97	0.47
2:H:110:THR:HG23	2:H:134:GLN:HG3	1.96	0.47
3:I:126:TYR:CZ	3:L:150:PRO:HA	2.49	0.47
3:C:86:ALA:HB1	1:D:52:GLU:HG2	1.97	0.46
3:L:57:ASN:O	3:L:64:PRO:HD3	2.15	0.46
3:F:169:GLU:OE1	3:F:172:LYS:HG2	2.16	0.46
1:G:117:PHE:CZ	1:J:214:PRO:HB3	2.50	0.46
1:J:26:CYS:SG	1:J:62:GLN:HA	2.56	0.46
1:J:63:LEU:HA	1:J:66:VAL:HG22	1.97	0.46
2:H:300:CYS:HA	2:H:301:PRO:HD3	1.77	0.46
3:C:98:ASP:OD1	3:C:98:ASP:N	2.47	0.46
1:D:225:TYR:O	1:D:229:LYS:HG2	2.16	0.46
1:G:110:THR:CG2	1:G:114:PRO:HD2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:176:LYS:HD3	2:H:203:ALA:HB3	1.98	0.46
1:J:302:PRO:HB2	1:J:305:GLN:HB2	1.97	0.46
1:A:141:GLN:OE1	1:A:180:ASP:N	2.45	0.46
1:G:89:MET:HA	1:G:92:TYR:CD2	2.50	0.46
2:H:117:ARG:NH2	2:H:312:ASP:OD1	2.42	0.46
3:C:105:ILE:HD11	3:C:211:LEU:HD11	1.97	0.46
1:D:43:MET:O	1:D:48:GLU:N	2.48	0.46
3:C:197:ILE:HD11	3:C:224:PRO:HD2	1.97	0.46
2:K:193:PRO:O	2:K:197:ARG:HG3	2.16	0.46
1:A:371:MET:HE2	1:A:391:ILE:HD12	1.98	0.46
1:D:31:ARG:NH1	7:F:402[A]:O2T:O7	2.48	0.46
1:A:152:ARG:NH1	1:A:190:PHE:O	2.50	0.45
3:I:115:ASP:OD2	3:I:141:TYR:OH	2.33	0.45
3:C:59:SER:HA	3:C:60:TYR:HA	1.62	0.45
1:D:154:GLY:HA3	1:D:163:GLN:NE2	2.32	0.45
2:H:94:SER:HA	2:H:311:LEU:HD21	1.98	0.45
2:K:90:LEU:HD13	2:K:95:LEU:HB2	1.98	0.45
2:B:11:ASN:ND2	2:B:72:ASP:OD1	2.44	0.45
1:D:134:GLN:HB3	1:D:299:ARG:HB2	1.97	0.45
1:D:196:ASN:OD1	1:D:197:HIS:N	2.45	0.45
1:G:205:PRO:HA	3:L:68:PRO:HG3	1.97	0.45
1:G:307:LYS:HA	1:G:342:LEU:HD23	1.97	0.45
3:C:29:PRO:HG2	4:C:401:SF4:S4	2.57	0.45
1:D:335:LEU:HD22	1:D:353:GLY:HA3	1.98	0.45
1:G:216:MET:C	1:G:219:PRO:HD2	2.42	0.45
1:J:17:VAL:HG13	1:J:22:VAL:HG21	1.98	0.45
3:C:47:ILE:HD11	3:C:84:ILE:HD12	1.99	0.45
3:L:229:TRP:HB2	3:L:231:PHE:CE2	2.52	0.45
1:D:79:GLY:HA2	1:D:104:MET:HG3	1.99	0.45
2:E:96:ILE:HD12	2:E:125:PHE:HB2	1.98	0.45
3:F:76:ASN:HB2	3:F:79:ALA:HB3	1.97	0.45
3:F:81:ALA:HB1	3:F:131:VAL:HG11	1.97	0.45
1:A:24:VAL:HG21	1:A:69:ALA:HB1	1.99	0.45
1:G:213:GLU:HB2	1:G:216:MET:HG2	1.99	0.45
1:G:282:ARG:NH2	2:H:75:GLU:OE1	2.48	0.45
2:H:244:MET:HE3	2:H:246:THR:HG22	1.99	0.45
1:J:79:GLY:HA2	1:J:104:MET:HG3	1.99	0.45
1:J:154:GLY:HA3	1:J:163:GLN:NE2	2.31	0.45
3:L:139:GLU:HB3	3:L:153:PRO:HD3	1.99	0.45
1:A:76:VAL:HG21	1:A:102:VAL:HG22	1.99	0.45
1:A:219:PRO:HB3	2:E:225:LEU:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:LYS:HA	1:A:342:LEU:HD23	1.99	0.45
2:B:186:LEU:HD22	2:B:196:MET:HE3	1.98	0.45
3:C:139:GLU:HB3	3:C:153:PRO:HD3	1.99	0.45
2:E:25:ASN:OD1	2:E:26:LYS:N	2.50	0.45
1:A:251:ASP:O	1:A:296:ARG:HD3	2.16	0.44
2:E:109:SER:OG	2:E:132:LEU:HA	2.17	0.44
2:E:244:MET:HE3	2:E:246:THR:HG22	1.99	0.44
1:J:273:MET:HG3	1:J:294:ARG:HG3	1.98	0.44
1:D:243:ALA:HB2	1:D:249:LYS:HA	1.99	0.44
2:E:94:SER:HA	2:E:311:LEU:HD21	1.99	0.44
1:G:218:PRO:HG2	1:J:361:GLU:HG3	1.98	0.44
3:I:59:SER:HA	3:I:60:TYR:HA	1.53	0.44
2:K:278:ARG:NH1	3:L:13:ASP:OD1	2.45	0.44
1:A:109:ARG:NH1	1:A:116:ASP:OD1	2.48	0.44
1:D:161:LEU:O	1:D:163:GLN:NE2	2.43	0.44
2:E:17:VAL:HG22	2:E:89:VAL:HB	1.99	0.44
2:B:93:GLU:O	2:B:96:ILE:HG13	2.18	0.44
1:D:213:GLU:HG2	1:D:215:ALA:H	1.82	0.44
3:F:197:ILE:HD11	3:F:224:PRO:HD2	2.00	0.44
1:G:33:TYR:HB3	1:G:81:SER:OG	2.17	0.44
1:J:279:VAL:HG21	1:J:368:PHE:HB2	2.00	0.44
1:D:252:PRO:HB2	1:D:253:TYR:CE2	2.53	0.44
1:D:261:ASP:O	1:D:289:LYS:HD3	2.18	0.44
2:H:110:THR:OG1	2:H:134:GLN:NE2	2.40	0.44
3:L:51:GLY:HA2	5:L:402:TPP:H71	1.98	0.44
1:D:371:MET:HE2	1:D:391:ILE:HD12	2.00	0.44
3:C:29:PRO:HB3	3:C:136:TYR:CD2	2.53	0.44
1:J:216:MET:C	1:J:219:PRO:HD2	2.43	0.44
2:K:295:LEU:HD22	3:L:237:ILE:HD12	1.99	0.44
3:L:248:MET:HG3	3:L:274:TYR:CD2	2.52	0.44
2:B:46:VAL:HG12	2:B:69:ILE:HG12	1.99	0.44
3:C:88:TYR:OH	1:D:200:LEU:O	2.29	0.44
2:E:31:ILE:O	2:E:35:LEU:HD13	2.18	0.44
2:E:110:THR:HG23	2:E:134:GLN:HB2	2.00	0.44
3:F:29:PRO:HG2	4:F:401:SF4:S4	2.58	0.44
1:D:307:LYS:HA	1:D:342:LEU:HD23	1.99	0.43
2:E:58:ARG:O	2:E:61:VAL:HG12	2.18	0.43
3:L:267:LYS:HD3	3:L:310:VAL:HG11	1.99	0.43
1:A:321:VAL:HG13	1:A:357:GLY:HA3	2.00	0.43
2:B:253:ARG:NH2	2:B:310:GLU:OE1	2.38	0.43
3:C:33:TYR:CE1	3:C:107:MET:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:154:GLY:HA3	1:G:163:GLN:NE2	2.33	0.43
3:I:58:THR:CG2	1:J:211:GLN:H	2.31	0.43
1:J:168:ASP:HB2	1:J:172:VAL:HG23	2.00	0.43
1:J:306:ILE:HG23	1:J:310:LEU:HD12	2.00	0.43
2:B:193:PRO:O	2:B:197:ARG:HG3	2.19	0.43
3:C:74:ILE:HD12	5:C:402:TPP:H72	2.00	0.43
3:F:138:ASN:HB2	3:F:225:CYS:HB2	2.01	0.43
1:G:65:VAL:O	3:L:83:GLY:HA3	2.19	0.43
1:G:152:ARG:HD2	1:G:190:PHE:O	2.18	0.43
3:I:2:LEU:HB2	3:I:212:ASN:HD21	1.84	0.43
1:A:78:THR:OG1	1:A:79:GLY:N	2.52	0.43
3:F:125:LEU:HB2	3:F:186:HIS:CD2	2.53	0.43
1:G:273:MET:O	1:G:277:LYS:HG3	2.18	0.43
1:G:284:ARG:HD2	1:G:284:ARG:HA	1.87	0.43
1:J:367:GLU:HG2	1:J:370:ARG:HH21	1.84	0.43
2:K:274:SER:HB2	3:L:19:PRO:HD3	2.01	0.43
1:A:26:CYS:SG	1:A:62:GLN:HA	2.58	0.43
3:F:229:TRP:HA	3:F:280:ARG:HD3	2.01	0.43
1:G:264:VAL:HG12	1:G:315:ALA:HB3	1.99	0.43
2:H:107:PRO:HA	2:H:131:ASN:HB3	2.00	0.43
3:I:229:TRP:HB2	3:I:231:PHE:CE2	2.54	0.43
3:L:229:TRP:HA	3:L:280:ARG:HD3	2.00	0.43
3:C:169:GLU:OE1	3:C:169:GLU:N	2.44	0.43
3:F:16:TYR:CD1	3:F:66:ARG:HD3	2.54	0.43
3:I:68:PRO:HG3	1:J:205:PRO:HA	2.00	0.43
3:I:75:THR:HG22	3:I:116:ILE:HG12	2.01	0.43
2:K:38:ALA:HB1	2:K:181:VAL:HG11	2.01	0.43
3:F:209:LYS:O	3:F:213:GLN:HG2	2.18	0.43
1:G:367:GLU:HG2	1:G:370:ARG:HH22	1.84	0.43
3:I:43:ASN:HA	1:J:202:PRO:HB3	2.00	0.43
1:D:251:ASP:HA	1:D:252:PRO:HD3	1.93	0.43
2:E:35:LEU:HD11	2:E:171:ALA:HB1	2.01	0.43
2:E:89:VAL:HG22	2:E:112:VAL:HB	2.00	0.43
3:I:125:LEU:HB2	3:I:186:HIS:CD2	2.53	0.43
1:A:261:ASP:O	1:A:289:LYS:HD3	2.19	0.42
2:B:120:ASP:OD1	2:B:206:ARG:NH1	2.51	0.42
3:F:224:PRO:HB2	3:F:236:THR:HG23	2.01	0.42
3:I:126:TYR:CE1	3:L:150:PRO:HA	2.54	0.42
1:J:109:ARG:HA	1:J:119:GLU:HA	2.01	0.42
3:C:225:CYS:HA	3:C:226:PRO:HD2	1.86	0.42
1:J:33:TYR:HB3	1:J:81:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:265:TYR:CD2	2:K:278:ARG:HD2	2.54	0.42
1:D:181:ILE:HA	1:D:182:PRO:HD3	1.83	0.42
2:E:107:PRO:HA	2:E:131:ASN:HB3	2.01	0.42
1:J:141:GLN:OE1	1:J:180:ASP:N	2.47	0.42
1:J:181:ILE:HA	1:J:182:PRO:HD3	1.81	0.42
2:E:278:ARG:NE	3:F:13:ASP:OD1	2.51	0.42
2:H:86:ASP:OD1	2:H:106:ARG:HD2	2.19	0.42
1:G:207:ILE:HD11	2:K:237:PRO:HD3	2.00	0.42
1:A:213:GLU:HG2	1:A:215:ALA:H	1.84	0.42
3:F:18:VAL:HG22	3:F:64:PRO:O	2.19	0.42
3:F:59:SER:HA	3:F:60:TYR:HA	1.61	0.42
3:F:180:LYS:HE3	3:F:259:TYR:CE2	2.54	0.42
1:A:365:HIS:NE2	2:B:75:GLU:O	2.48	0.42
3:C:63:GLY:HA3	1:D:207:ILE:HD13	2.01	0.42
3:I:57:ASN:O	3:I:64:PRO:HD3	2.20	0.42
1:A:128:ARG:HB3	1:A:299:ARG:HH21	1.84	0.42
2:B:59:ILE:HB	3:C:25:ALA:H	1.84	0.42
1:D:89:MET:HA	1:D:92:TYR:CE2	2.54	0.42
3:I:63:GLY:HA3	1:J:207:ILE:HD13	2.01	0.42
3:L:104:ILE:HB	3:L:131:VAL:HG12	2.01	0.42
3:L:298:ALA:O	3:L:302:GLU:HG3	2.20	0.42
1:A:17:VAL:HG13	1:A:22:VAL:HG21	2.01	0.42
1:A:306:ILE:HG23	1:A:310:LEU:HD12	2.02	0.42
3:C:264:LYS:HD2	3:C:264:LYS:N	2.34	0.42
1:D:93:SER:HB3	1:D:94:PRO:HD3	2.00	0.42
2:E:246:THR:OG1	2:E:289:TYR:O	2.37	0.42
3:F:33:TYR:CZ	3:F:107:MET:HG3	2.55	0.42
3:I:128:GLY:HA2	3:I:216:PRO:HG3	2.01	0.42
2:K:58:ARG:O	2:K:61:VAL:HG12	2.20	0.42
2:B:209:PRO:HA	2:B:210:PRO:HD3	1.96	0.41
3:F:264:LYS:HA	3:F:264:LYS:HD2	1.83	0.41
1:G:52:GLU:HG2	3:L:90:ALA:HB2	2.01	0.41
2:K:291:LYS:NZ	3:L:23:THR:O	2.44	0.41
1:D:367:GLU:O	1:D:371:MET:HG3	2.20	0.41
2:E:244:MET:HE2	3:F:22:ARG:HG3	2.02	0.41
3:I:2:LEU:HB3	3:I:208:ARG:HH11	1.85	0.41
3:F:57:ASN:O	3:F:64:PRO:HD3	2.20	0.41
1:G:43:MET:O	1:G:48:GLU:N	2.52	0.41
1:G:128:ARG:HG2	1:G:298:PHE:CE2	2.55	0.41
1:G:302:PRO:HB2	1:G:305:GLN:HB2	2.02	0.41
3:I:29:PRO:HB3	3:I:136:TYR:CD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:286:LEU:HD23	1:J:286:LEU:HA	1.91	0.41
3:L:179:PRO:HG3	3:L:220:HIS:CB	2.50	0.41
1:A:18:ARG:HB2	1:A:43:MET:HE1	2.02	0.41
1:G:76:VAL:HG21	1:G:102:VAL:HG22	2.02	0.41
1:G:214:PRO:HB3	1:J:117:PHE:CE2	2.55	0.41
2:H:18:TRP:CH2	2:H:64:ARG:HD3	2.56	0.41
2:H:93:GLU:O	2:H:96:ILE:HG12	2.20	0.41
3:I:138:ASN:HB2	3:I:225:CYS:HB2	2.02	0.41
1:J:82:GLY:HA3	1:J:109:ARG:HD2	2.03	0.41
1:A:18:ARG:HB2	1:A:43:MET:CE	2.50	0.41
2:B:86:ASP:OD2	2:B:106:ARG:NH1	2.36	0.41
3:C:169:GLU:OE2	3:C:172:LYS:HG2	2.21	0.41
2:E:271:CYS:HA	2:E:272:PRO:HD3	1.91	0.41
3:I:33:TYR:CZ	3:I:107:MET:HG3	2.56	0.41
2:K:96:ILE:HD13	2:K:122:ILE:HD12	2.02	0.41
2:E:120:ASP:O	2:E:124:GLU:HG3	2.20	0.41
1:G:52:GLU:HG3	3:L:86:ALA:O	2.21	0.41
1:J:89:MET:HA	1:J:92:TYR:CE2	2.56	0.41
1:A:108:ASP:HB3	1:A:120:GLU:O	2.20	0.41
2:B:57:ASP:OD1	2:B:57:ASP:N	2.46	0.41
1:G:181:ILE:HA	1:G:182:PRO:HD3	1.81	0.41
2:K:271:CYS:HA	2:K:272:PRO:HD3	1.95	0.41
3:L:3:ASP:O	3:L:208:ARG:NH1	2.41	0.41
3:L:222:HIS:HB2	3:L:249:PHE:CE1	2.55	0.41
1:A:225:TYR:O	1:A:229:LYS:HG2	2.21	0.41
2:B:17:VAL:HG22	2:B:89:VAL:HB	2.02	0.41
3:C:152:THR:HB	3:C:171:LYS:HE3	2.02	0.41
3:I:163:PRO:HB3	3:I:283:HIS:CD2	2.56	0.41
3:I:232:PRO:HB2	3:I:235:LYS:HG2	2.03	0.41
1:J:391:ILE:HA	1:J:392:PRO:HA	1.86	0.41
2:K:129:THR:HB	2:K:132:LEU:HB3	2.02	0.41
1:A:112:ASP:OD1	1:A:115:GLY:HA2	2.20	0.41
2:B:170:ILE:O	2:B:174:VAL:HG23	2.20	0.41
2:B:279:THR:OG1	2:B:282:GLY:O	2.35	0.41
3:C:125:LEU:HB2	3:C:186:HIS:CD2	2.56	0.41
1:D:273:MET:O	1:D:277:LYS:HG3	2.21	0.41
3:F:169:GLU:OE2	3:F:169:GLU:N	2.51	0.41
1:G:37:MET:HE1	1:G:78:THR:HA	2.03	0.41
1:J:32:PRO:HD3	1:J:109:ARG:HB2	2.02	0.41
2:K:192:ASN:HA	2:K:193:PRO:HD2	1.97	0.41
1:A:335:LEU:HD22	1:A:353:GLY:HA3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:115:THR:HG21	2:H:122:ILE:HD11	2.03	0.41
3:I:152:THR:HB	3:I:171:LYS:HE3	2.03	0.41
3:C:18:VAL:HG22	3:C:64:PRO:O	2.21	0.40
2:E:209:PRO:HA	2:E:210:PRO:HD3	1.92	0.40
1:G:17:VAL:HG13	1:G:22:VAL:HG21	2.02	0.40
1:G:343:TYR:CZ	1:J:346:GLY:HA2	2.56	0.40
2:H:116:LYS:HE2	2:H:116:LYS:HB3	1.82	0.40
1:J:367:GLU:HG2	1:J:370:ARG:NH2	2.36	0.40
1:A:89:MET:HA	1:A:92:TYR:CE2	2.56	0.40
1:D:301:PHE:HA	1:D:302:PRO:HD3	1.88	0.40
1:G:142:GLU:OE2	1:G:294:ARG:NH1	2.53	0.40
1:G:200:LEU:HD21	3:L:70:ILE:HB	2.03	0.40
3:I:54:TYR:O	3:I:58:THR:HG22	2.21	0.40
3:L:222:HIS:CE1	3:L:229:TRP:HH2	2.39	0.40
3:L:259:TYR:HB3	3:L:303:VAL:HG13	2.02	0.40
3:F:125:LEU:HD23	3:F:216:PRO:HB2	2.02	0.40
3:I:68:PRO:HD3	1:J:205:PRO:HB3	2.04	0.40
2:B:18:TRP:CH2	2:B:64:ARG:HD3	2.56	0.40
3:C:227:LYS:HD3	4:C:401:SF4:S2	2.62	0.40
3:I:29:PRO:HG2	4:I:401:SF4:S4	2.62	0.40
2:K:68:ARG:HH22	2:K:79:ILE:HB	1.86	0.40
2:K:293:CYS:O	3:L:234:ASP:HB3	2.22	0.40
3:L:169:GLU:OE1	3:L:169:GLU:N	2.53	0.40
1:A:33:TYR:HB3	1:A:81:SER:OG	2.21	0.40
2:E:115:THR:HG21	2:E:122:ILE:HD11	2.03	0.40
3:F:180:LYS:HE3	3:F:259:TYR:HE2	1.87	0.40
1:G:203:ARG:NH2	3:L:99:ALA:HB2	2.37	0.40
1:G:225:TYR:CE1	1:G:229:LYS:HD3	2.56	0.40
1:G:290:VAL:HB	1:G:372:PHE:CE1	2.57	0.40
1:J:251:ASP:O	1:J:296:ARG:HD3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	392/395 (99%)	378 (96%)	14 (4%)	0	100	100
1	D	392/395 (99%)	379 (97%)	12 (3%)	1 (0%)	36	55
1	G	392/395 (99%)	378 (96%)	14 (4%)	0	100	100
1	J	392/395 (99%)	378 (96%)	13 (3%)	1 (0%)	36	55
2	B	280/315 (89%)	272 (97%)	8 (3%)	0	100	100
2	E	289/315 (92%)	280 (97%)	9 (3%)	0	100	100
2	H	287/315 (91%)	277 (96%)	10 (4%)	0	100	100
2	K	274/315 (87%)	267 (97%)	7 (3%)	0	100	100
3	C	310/314 (99%)	301 (97%)	8 (3%)	1 (0%)	36	55
3	F	308/314 (98%)	299 (97%)	8 (3%)	1 (0%)	36	55
3	I	307/314 (98%)	296 (96%)	10 (3%)	1 (0%)	36	55
3	L	308/314 (98%)	298 (97%)	9 (3%)	1 (0%)	36	55
All	All	3931/4096 (96%)	3803 (97%)	122 (3%)	6 (0%)	43	63

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	L	146	ILE
3	F	146	ILE
3	I	146	ILE
3	C	146	ILE
1	D	114	PRO
1	J	114	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	319/322 (99%)	312 (98%)	7 (2%)	45	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	317/322 (98%)	313 (99%)	4 (1%)	61	82
1	G	318/322 (99%)	312 (98%)	6 (2%)	50	76
1	J	318/322 (99%)	314 (99%)	4 (1%)	61	82
2	B	233/256 (91%)	233 (100%)	0	100	100
2	E	227/256 (89%)	227 (100%)	0	100	100
2	H	236/256 (92%)	236 (100%)	0	100	100
2	K	219/256 (86%)	219 (100%)	0	100	100
3	C	249/250 (100%)	248 (100%)	1 (0%)	84	93
3	F	245/250 (98%)	245 (100%)	0	100	100
3	I	244/250 (98%)	243 (100%)	1 (0%)	84	93
3	L	246/250 (98%)	243 (99%)	3 (1%)	63	83
All	All	3171/3312 (96%)	3145 (99%)	26 (1%)	73	88

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	33	TYR
1	A	119	GLU
1	A	211	GLN
1	A	292	VAL
1	A	333	VAL
1	A	362	VAL
3	C	33	TYR
1	D	33	TYR
1	D	119	GLU
1	D	333	VAL
1	D	362	VAL
1	G	33	TYR
1	G	36	ILE
1	G	110	THR
1	G	119	GLU
1	G	333	VAL
1	G	363	VAL
3	I	94	LYS
1	J	33	TYR
1	J	36	ILE
1	J	119	GLU

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Mol	Chain	Res	Type
1	J	260	ASP
3	L	33	TYR
3	L	251	LEU
3	L	266	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	272	HIS
1	A	285	ASN
1	A	324	ASN
3	C	43	ASN
3	C	178	ASN
1	D	6	ASN
1	D	226	GLN
1	D	324	ASN
2	E	11	ASN
2	E	131	ASN
3	F	230	GLN
1	G	6	ASN
1	G	130	GLN
1	G	197	HIS
1	G	269	GLN
1	G	285	ASN
2	H	134	GLN
2	H	207	GLN
3	I	222	HIS
1	J	130	GLN
1	J	226	GLN
1	J	272	HIS
2	K	204	GLN
3	L	213	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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
5	TPP	L	402	6	26,27,27	1.88	7 (26%)	38,40,40	1.89	10 (26%)
4	SF4	F	401	3	0,12,12	-	-	-		
4	SF4	L	401	3	0,12,12	-	-	-		
4	SF4	K	401	2	0,12,12	-	-	-		
4	SF4	C	401	3	0,12,12	-	-	-		
4	SF4	I	401	3	0,12,12	-	-	-		
4	SF4	E	401	2	0,12,12	-	-	-		
5	TPP	C	402	6	26,27,27	1.87	7 (26%)	38,40,40	1.83	9 (23%)
4	SF4	E	402	2	0,12,12	-	-	-		
5	TPP	I	402	6	26,27,27	1.90	8 (30%)	38,40,40	1.79	9 (23%)
4	SF4	B	401	2	0,12,12	-	-	-		
4	SF4	K	402	2	0,12,12	-	-	-		
4	SF4	H	401	2	0,12,12	-	-	-		
4	SF4	H	402	2	0,12,12	-	-	-		
4	SF4	B	402	2	0,12,12	-	-	-		

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
5	TPP	L	402	6	-	1/17/17/17	0/2/2/2
4	SF4	F	401	3	-	-	0/6/5/5
4	SF4	L	401	3	-	-	0/6/5/5
4	SF4	K	401	2	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SF4	C	401	3	-	-	0/6/5/5
4	SF4	I	401	3	-	-	0/6/5/5
4	SF4	E	401	2	-	-	0/6/5/5
4	SF4	H	402	2	-	-	0/6/5/5
4	SF4	K	402	2	-	-	0/6/5/5
4	SF4	E	402	2	-	-	0/6/5/5
4	SF4	B	401	2	-	-	0/6/5/5
5	TPP	C	402	6	-	1/17/17/17	0/2/2/2
4	SF4	H	401	2	-	-	0/6/5/5
5	TPP	I	402	6	-	4/17/17/17	0/2/2/2
4	SF4	B	402	2	-	-	0/6/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	402	TPP	C7'-C5'	3.93	1.58	1.51
5	I	402	TPP	C5-S1	-3.71	1.63	1.72
5	C	402	TPP	C5-S1	-3.70	1.63	1.72
5	C	402	TPP	C7'-C5'	3.66	1.57	1.51
5	I	402	TPP	C4'-N4'	3.63	1.43	1.34
5	L	402	TPP	C5-S1	-3.62	1.63	1.72
5	I	402	TPP	C7'-C5'	3.59	1.57	1.51
5	L	402	TPP	C4'-N4'	3.58	1.43	1.34
5	C	402	TPP	C4'-N4'	3.53	1.43	1.34
5	C	402	TPP	C2-N3	3.09	1.39	1.32
5	L	402	TPP	C2-N3	3.05	1.39	1.32
5	I	402	TPP	C2-N3	2.99	1.39	1.32
5	I	402	TPP	C5-C4	2.79	1.41	1.35
5	I	402	TPP	C7'-N3	-2.75	1.42	1.48
5	C	402	TPP	C7'-N3	-2.74	1.42	1.48
5	L	402	TPP	C7'-N3	-2.61	1.42	1.48
5	L	402	TPP	C5-C4	2.54	1.40	1.35
5	C	402	TPP	C5-C4	2.51	1.40	1.35
5	L	402	TPP	C6'-C5'	2.40	1.42	1.37
5	C	402	TPP	C6'-C5'	2.36	1.42	1.37
5	I	402	TPP	C6'-C5'	2.30	1.42	1.37
5	I	402	TPP	C4-N3	-2.04	1.35	1.39

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	402	TPP	C2-S1-C5	-4.69	88.11	91.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	402	TPP	C2-N3-C4	-4.38	108.14	114.06
5	C	402	TPP	C2-N3-C4	-4.33	108.20	114.06
5	C	402	TPP	C2-S1-C5	-4.28	88.38	91.22
5	L	402	TPP	S1-C2-N3	4.19	117.61	112.30
5	I	402	TPP	C2-N3-C4	-4.02	108.63	114.06
5	I	402	TPP	C2-S1-C5	-3.94	88.60	91.22
5	C	402	TPP	S1-C2-N3	3.92	117.26	112.30
5	I	402	TPP	S1-C2-N3	3.83	117.15	112.30
5	L	402	TPP	C4-C5-S1	3.64	116.25	110.56
5	C	402	TPP	C4-C5-S1	3.51	116.03	110.56
5	I	402	TPP	C4-C5-S1	3.39	115.84	110.56
5	C	402	TPP	C6'-N1'-C2'	3.18	121.30	116.07
5	L	402	TPP	C6'-N1'-C2'	3.17	121.28	116.07
5	I	402	TPP	C6'-N1'-C2'	3.07	121.11	116.07
5	I	402	TPP	C5'-C6'-N1'	-2.65	119.52	123.83
5	C	402	TPP	C5'-C6'-N1'	-2.64	119.54	123.83
5	L	402	TPP	C5'-C6'-N1'	-2.60	119.61	123.83
5	I	402	TPP	N4'-C4'-N3'	2.44	120.31	117.03
5	C	402	TPP	N4'-C4'-N3'	2.36	120.22	117.03
5	C	402	TPP	C5'-C7'-N3	2.35	118.03	112.97
5	L	402	TPP	N4'-C4'-N3'	2.27	120.08	117.03
5	C	402	TPP	N1'-C2'-N3'	-2.22	121.83	125.53
5	L	402	TPP	N1'-C2'-N3'	-2.21	121.86	125.53
5	I	402	TPP	C5'-C7'-N3	2.19	117.69	112.97
5	I	402	TPP	N1'-C2'-N3'	-2.14	121.97	125.53
5	L	402	TPP	CM4-C4-N3	2.12	125.47	120.57
5	L	402	TPP	C5'-C7'-N3	2.12	117.54	112.97

There are no chirality outliers.

All (6) torsion outliers are listed below:

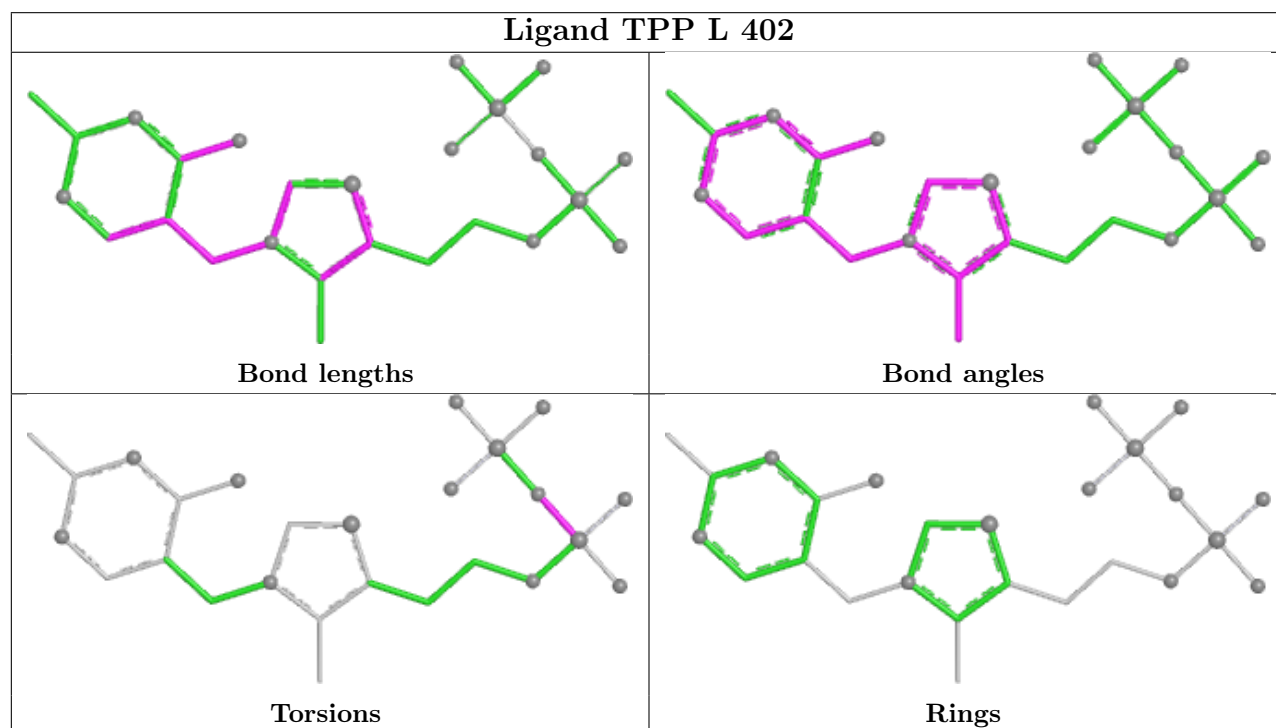
Mol	Chain	Res	Type	Atoms
5	L	402	TPP	PB-O3A-PA-O7
5	I	402	TPP	C4'-C5'-C7'-N3
5	I	402	TPP	C6'-C5'-C7'-N3
5	C	402	TPP	C5-C6-C7-O7
5	I	402	TPP	C5-C6-C7-O7
5	I	402	TPP	S1-C5-C6-C7

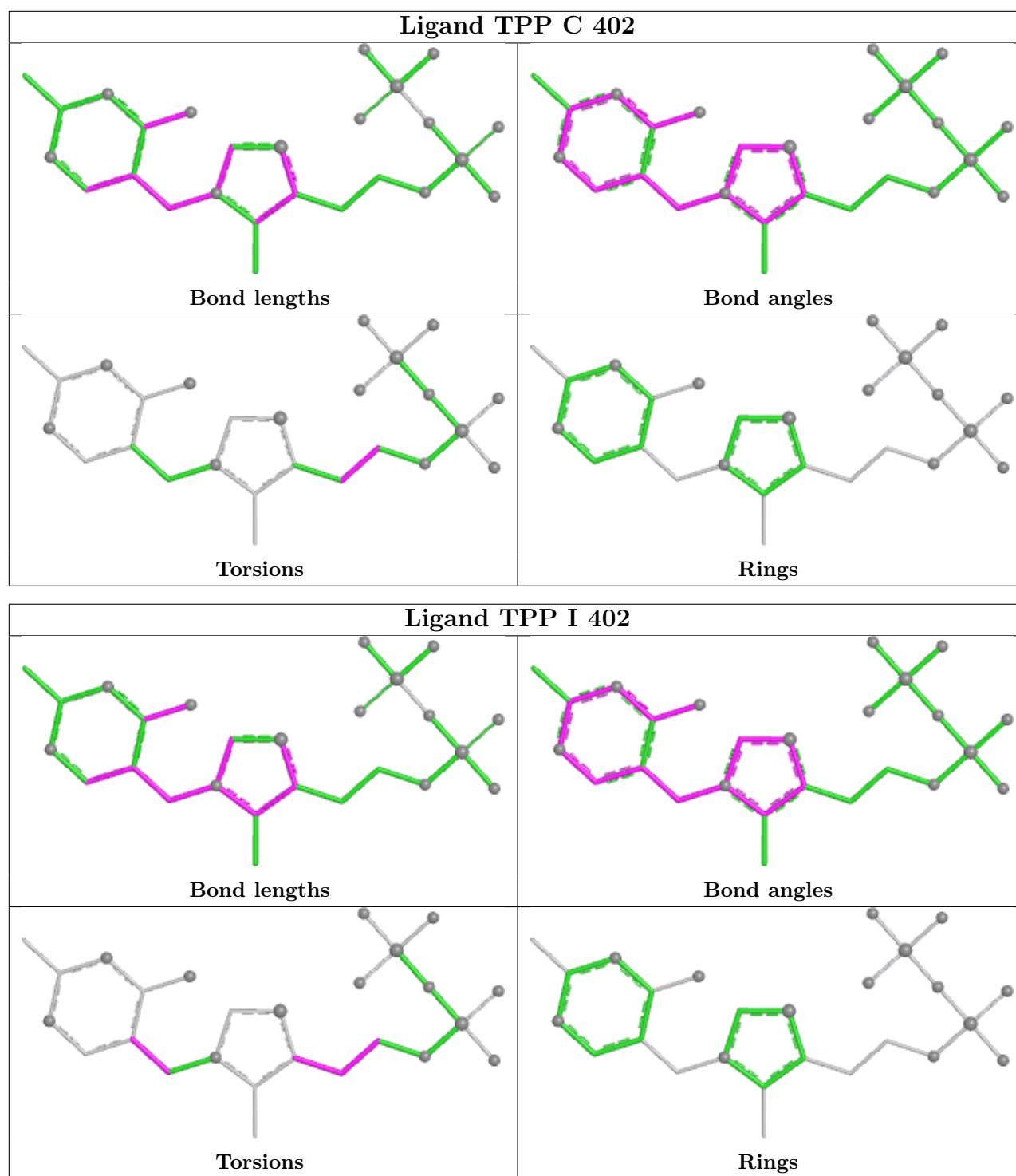
There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	402	TPP	1	0
4	F	401	SF4	2	0
4	L	401	SF4	1	0
4	C	401	SF4	2	0
4	I	401	SF4	1	0
5	C	402	TPP	1	0
5	I	402	TPP	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	394/395 (99%)	-1.58	0 100 100	21, 32, 47, 55	0
1	D	394/395 (99%)	-1.57	0 100 100	22, 35, 47, 52	0
1	G	394/395 (99%)	-1.65	0 100 100	19, 25, 33, 44	0
1	J	394/395 (99%)	-1.59	0 100 100	21, 30, 38, 44	0
2	B	286/315 (90%)	-1.50	0 100 100	24, 42, 58, 69	0
2	E	295/315 (93%)	-1.10	0 100 100	26, 66, 105, 117	0
2	H	291/315 (92%)	-1.49	0 100 100	27, 40, 56, 61	0
2	K	280/315 (88%)	-1.40	0 100 100	26, 54, 76, 82	0
3	C	312/314 (99%)	-1.63	0 100 100	20, 25, 38, 51	0
3	F	310/314 (98%)	-1.58	0 100 100	20, 29, 41, 54	0
3	I	309/314 (98%)	-1.56	0 100 100	25, 33, 45, 62	0
3	L	310/314 (98%)	-1.53	0 100 100	25, 35, 47, 60	0
All	All	3969/4096 (96%)	-1.52	0 100 100	19, 33, 67, 117	0

There are no RSRZ outliers to report.

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

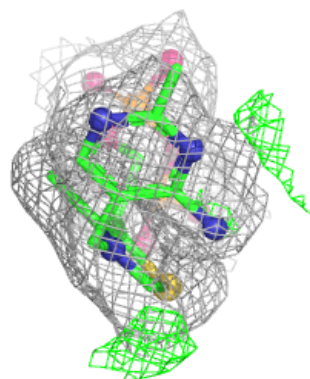
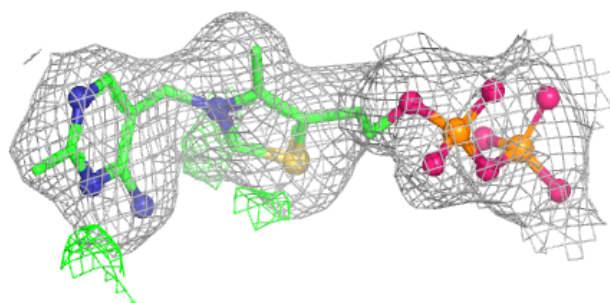
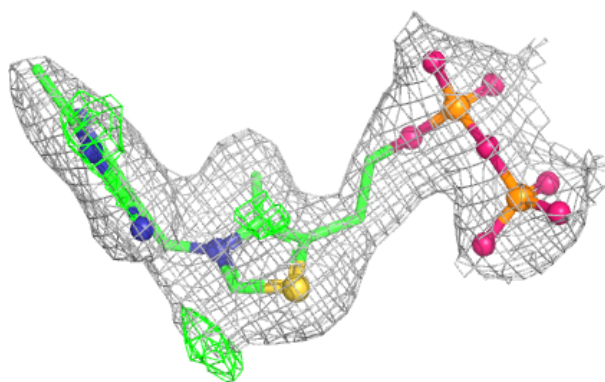
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	TPP	I	402	26/26	0.99	0.03	22,26,27,28	0
5	TPP	L	402	26/26	0.99	0.03	25,31,33,33	0
4	SF4	C	401	8/8	1.00	0.01	21,23,24,25	0
4	SF4	E	401	8/8	1.00	0.01	27,32,34,34	0
4	SF4	E	402	8/8	1.00	0.01	30,32,33,35	0
4	SF4	F	401	8/8	1.00	0.01	24,30,31,31	0
4	SF4	H	401	8/8	1.00	0.01	26,30,34,35	0
4	SF4	H	402	8/8	1.00	0.02	31,34,35,36	0
4	SF4	I	401	8/8	1.00	0.01	25,28,31,32	0
4	SF4	K	401	8/8	1.00	0.01	33,35,37,37	0
4	SF4	K	402	8/8	1.00	0.01	31,33,35,37	0
4	SF4	L	401	8/8	1.00	0.01	31,31,33,36	0
5	TPP	C	402	26/26	1.00	0.02	19,22,23,24	0
4	SF4	B	401	8/8	1.00	0.01	23,25,28,28	0
4	SF4	B	402	8/8	1.00	0.01	28,29,33,34	0
6	MG	C	403	1/1	1.00	0.01	19,19,19,19	0
6	MG	F	403	1/1	1.00	0.02	26,26,26,26	0
6	MG	I	403	1/1	1.00	0.00	26,26,26,26	0
6	MG	L	403	1/1	1.00	0.01	32,32,32,32	0
7	O2T	F	402[A]	32/32	1.00	0.03	24,28,31,31	6

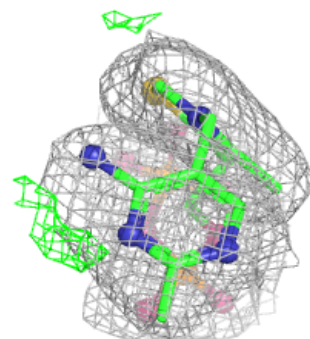
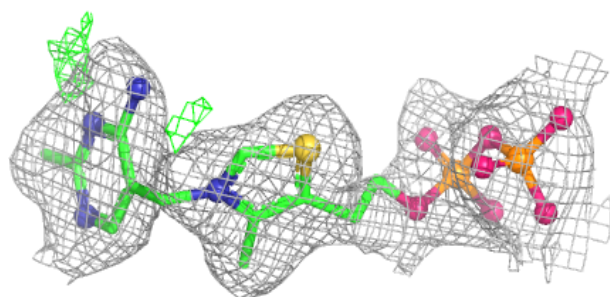
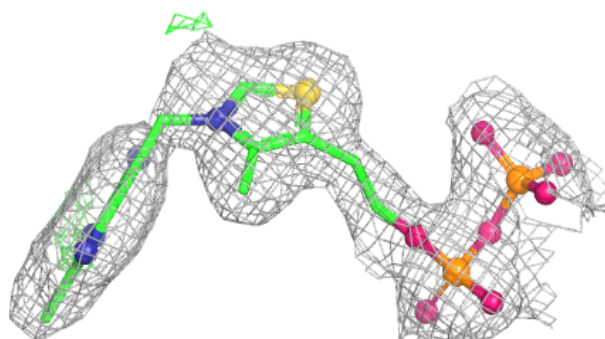
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

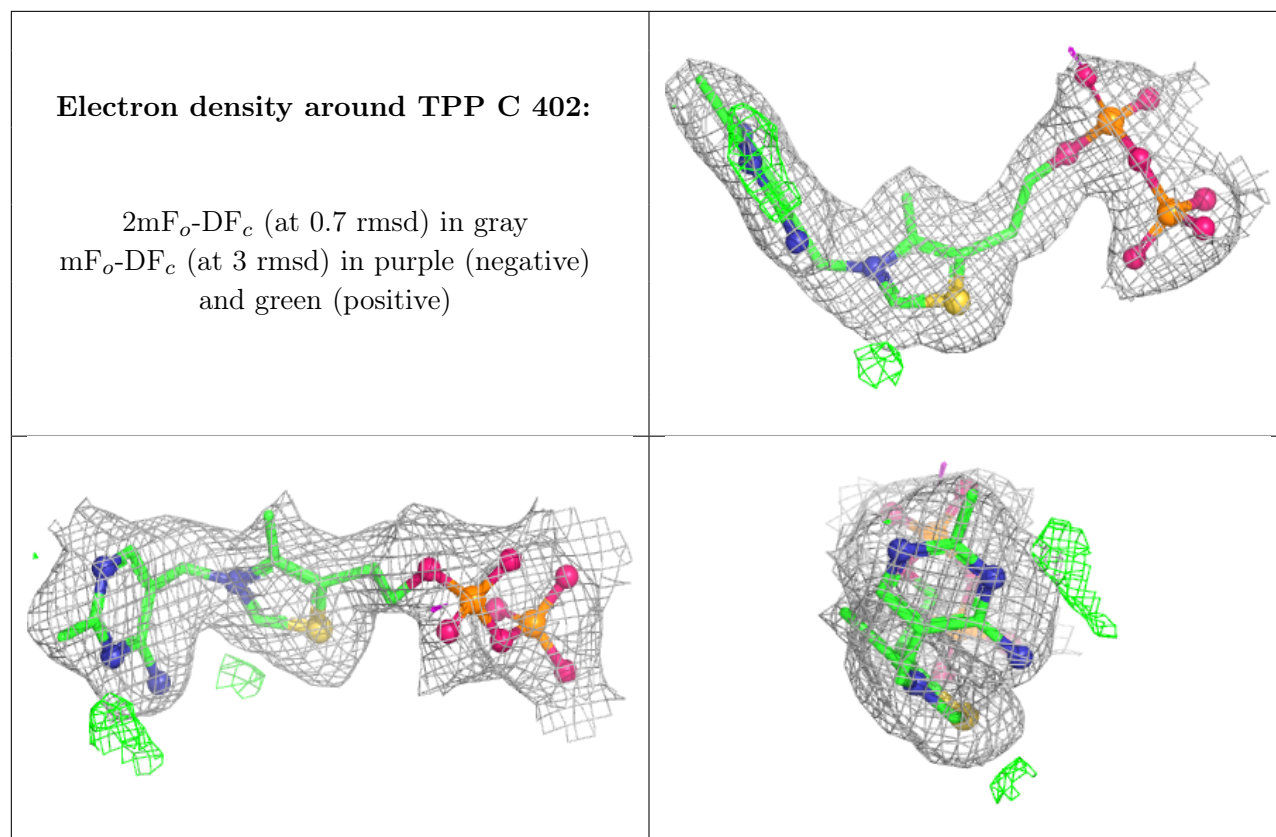
Electron density around TPP I 402:

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

**Electron density around TPP L 402:**

$2mF_o-DF_c$ (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.