



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:12 AM UTC

PDB ID : 1W88 / pdb_00001w88
Title : The crystal structure of pyruvate dehydrogenase E1(D180N,E183Q) bound to the peripheral subunit binding domain of E2
Authors : Frank, R.A.W.; Pratap, J.V.; Pei, X.Y.; Perham, R.N.; Luisi, B.F.
Deposited on : 2004-09-16
Resolution : 2.30 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

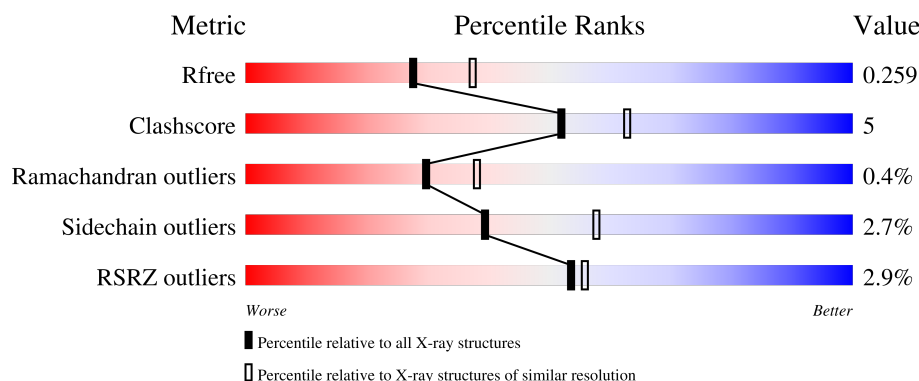
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	368	% 82% 11% 7%
1	C	368	2% 83% 9% 7%
1	E	368	3% 75% 12% 11%
1	G	368	2% 84% 5% 10%
2	B	324	% 86% 12%

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Mol	Chain	Length	Quality of chain
2	D	324	% 90% 9% .
2	F	324	2% 84% 10% . .
2	H	324	3% 88% 9% . .
3	I	49	24% 63% 12% 6% 18%
3	J	49	33% 63% 18% 6% 12%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 21934 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 PYRUVATE DEHYDROGENASE E1 COMPONENT, ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	1
			2592	1666	444	475	7			
1	C	343	Total	C	N	O	S	0	0	0
			2659	1705	451	496	7			
1	E	328	Total	C	N	O	S	0	1	0
			2521	1614	431	469	7			
1	G	331	Total	C	N	O	S	0	1	0
			2565	1645	434	479	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	180	ASN	ASP	engineered mutation	UNP P21873
A	183	GLN	GLU	engineered mutation	UNP P21873
C	180	ASN	ASP	engineered mutation	UNP P21873
C	183	GLN	GLU	engineered mutation	UNP P21873
E	180	ASN	ASP	engineered mutation	UNP P21873
E	183	GLN	GLU	engineered mutation	UNP P21873
G	180	ASN	ASP	engineered mutation	UNP P21873
G	183	GLN	GLU	engineered mutation	UNP P21873

- Molecule 2 is a protein called PYRUVATE DEHYDROGENASE E1 COMPONENT, BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	324	Total	C	N	O	S	0	1	0
			2428	1546	417	457	8			
2	D	324	Total	C	N	O	S	0	0	0
			2489	1586	424	471	8			
2	F	310	Total	C	N	O	S	0	1	0
			2369	1507	407	447	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	317	Total	C	N	O	S	0	0	0
			2416	1537	411	460	8			

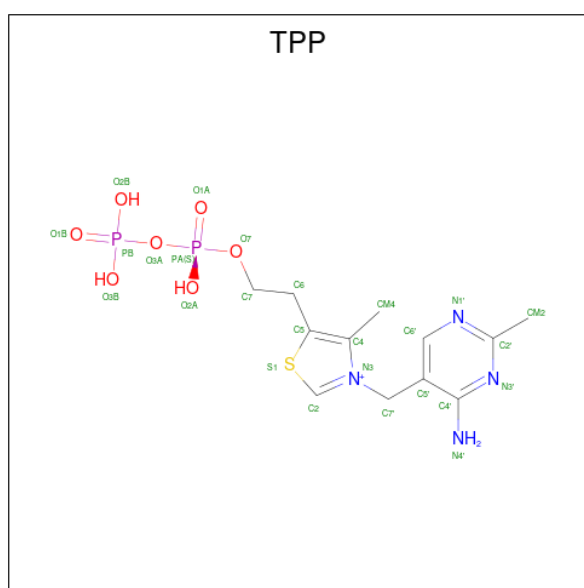
- Molecule 3 is a protein called DIHYDROLIPOYLLYSINE-RESIDUE ACETYLTRANSFERASE COMPONENT OF PYRUVATE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	40	Total	C	N	O	S	0	0	0
			303	192	55	55	1			
3	J	43	Total	C	N	O	S	0	0	1
			311	195	59	56	1			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	E	1	Total	Mg	0	0
			1	1		
4	G	1	Total	Mg	0	0
			1	1		

- Molecule 5 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: C₁₂H₁₉N₄O₇P₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	
5	C	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	
5	E	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	
5	G	1	Total	C	N	O	P	S	
			26	12	4	7	2	1	

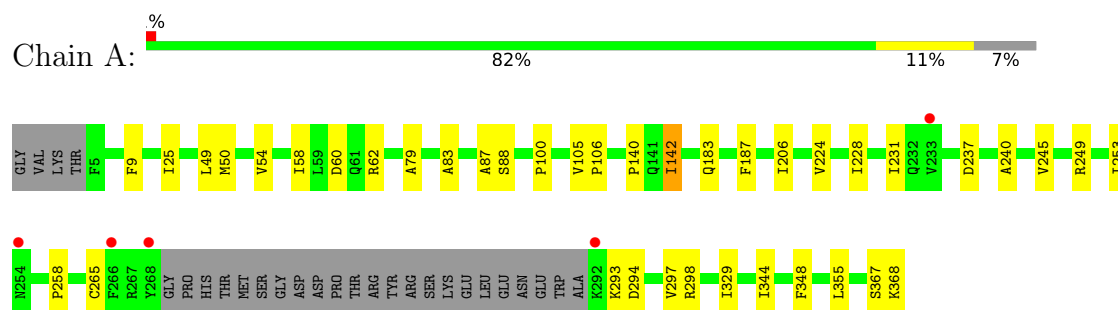
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	96	Total	O		
			96	96	0	0
6	B	124	Total	O		
			124	124	0	0
6	C	168	Total	O		
			168	168	0	0
6	D	209	Total	O		
			209	209	0	0
6	E	78	Total	O		
			78	78	0	0
6	F	194	Total	O		
			194	194	0	0
6	G	88	Total	O		
			88	88	0	0
6	H	161	Total	O		
			161	161	0	0
6	I	24	Total	O		
			24	24	0	0
6	J	31	Total	O		
			31	31	0	0

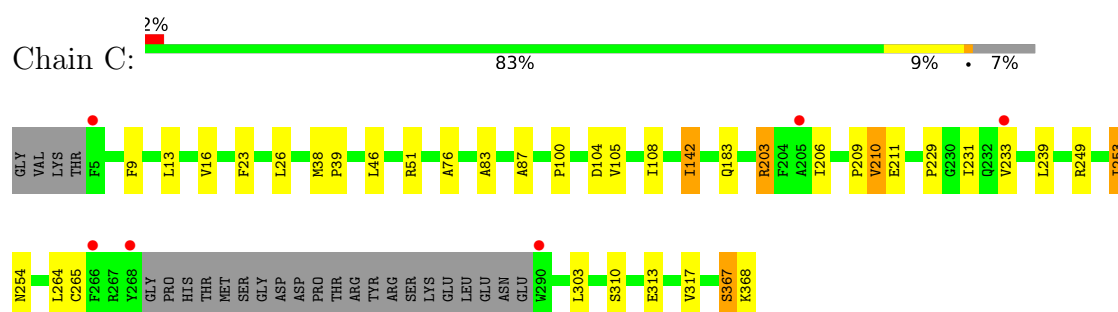
3 Residue-property plots [i](#)

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.

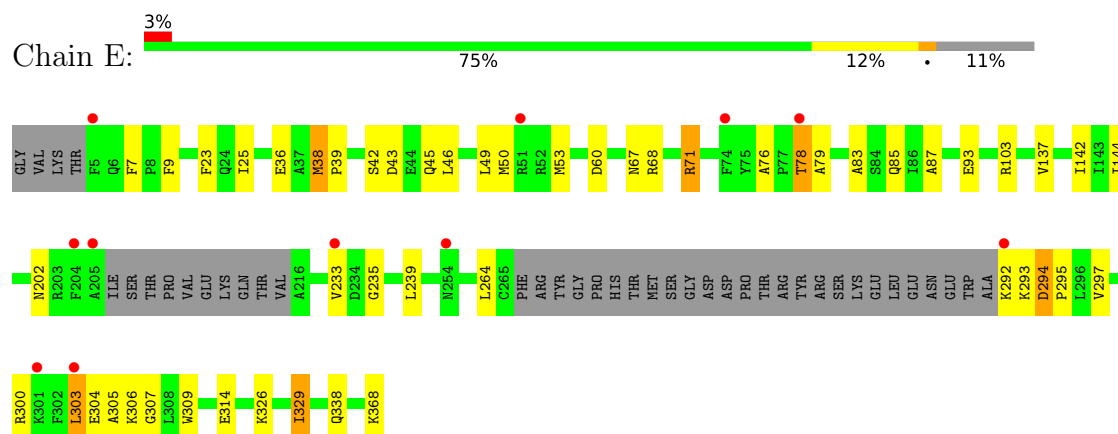
- Molecule 1: PYRUVATE DEHYDROGENASE E1 COMPONENT, ALPHA SUBUNIT



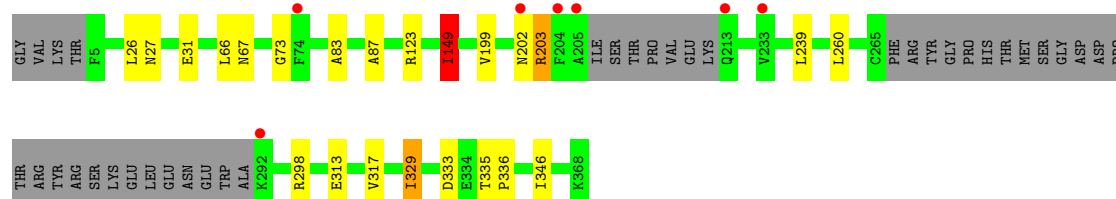
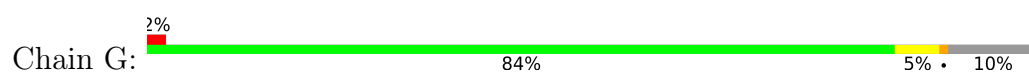
- Molecule 1: PYRUVATE DEHYDROGENASE E1 COMPONENT, ALPHA SUBUNIT



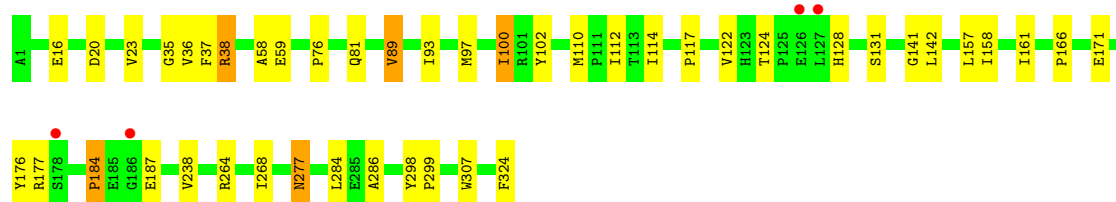
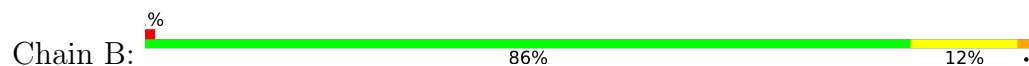
- Molecule 1: PYRUVATE DEHYDROGENASE E1 COMPONENT, ALPHA SUBUNIT



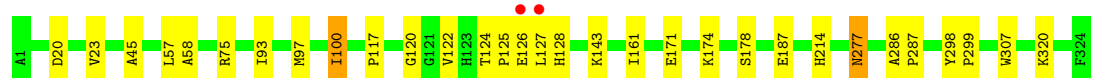
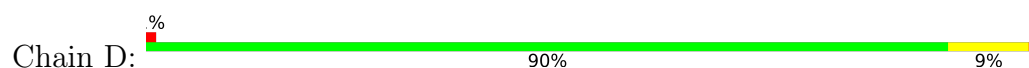
- Molecule 1: PYRUVATE DEHYDROGENASE E1 COMPONENT, ALPHA SUBUNIT



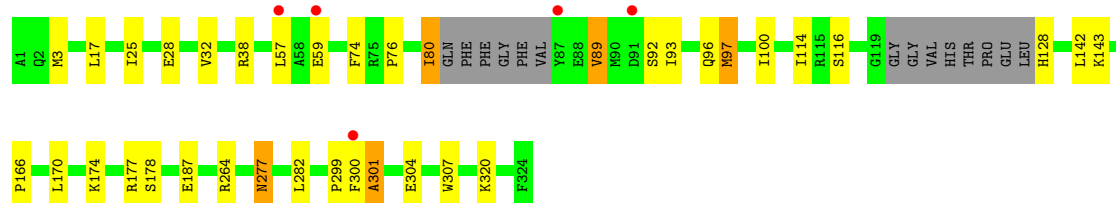
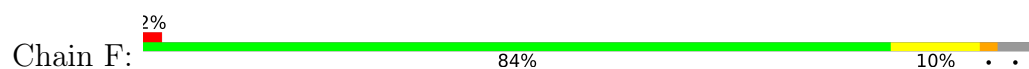
- Molecule 2: PYRUVATE DEHYDROGENASE E1 COMPONENT, BETA SUBUNIT



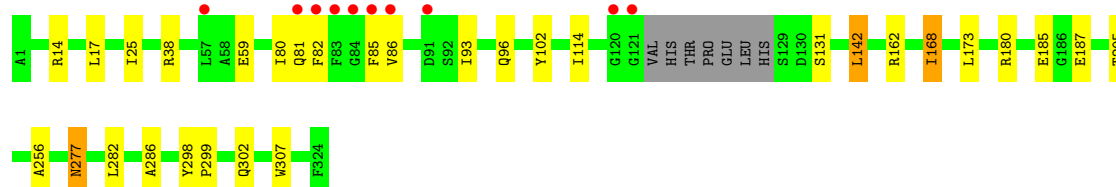
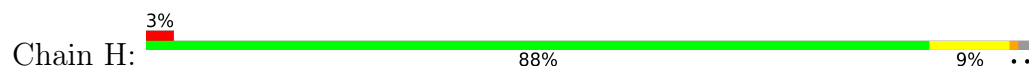
- Molecule 2: PYRUVATE DEHYDROGENASE E1 COMPONENT, BETA SUBUNIT



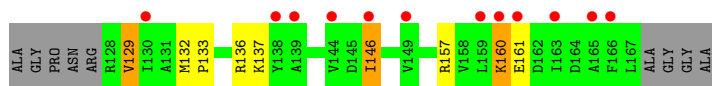
- Molecule 2: PYRUVATE DEHYDROGENASE E1 COMPONENT, BETA SUBUNIT



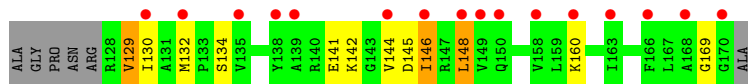
- Molecule 2: PYRUVATE DEHYDROGENASE E1 COMPONENT, BETA SUBUNIT



● Molecule 3: DIHYDROLIPOYLLYSINE-RESIDUE ACETYLTRANSFERASE COMPONENT OF PYRUVATE



● Molecule 3: DIHYDROLIPOYLLYSINE-RESIDUE ACETYLTRANSFERASE COMPONENT OF PYRUVATE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.18Å 133.69Å 245.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.30 19.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.98-2.30) 99.8 (19.98-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.199 , 0.262 0.199 , 0.259	Depositor DCC
R_{free} test set	6706 reflections (3.80%)	wwPDB-VP
Wilson B-factor (Å ²)	33.5	Xtriage
Anisotropy	0.564	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21934	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 7.15% 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, 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	0.65	1/2647 (0.0%)	0.89	0/3592
1	C	0.67	1/2715 (0.0%)	0.89	4/3677 (0.1%)
1	E	0.59	0/2571	0.87	0/3482
1	G	0.58	0/2617	0.89	1/3541 (0.0%)
2	B	0.72	0/2476	0.91	0/3369
2	D	0.71	0/2535	0.91	0/3437
2	F	0.72	0/2411	0.91	1/3267 (0.0%)
2	H	0.69	0/2458	0.90	1/3330 (0.0%)
3	I	0.62	0/305	1.01	0/408
3	J	0.76	1/313 (0.3%)	0.92	0/420
All	All	0.67	3/21048 (0.0%)	0.90	7/28523 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	E	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	169	GLY	C-N	-6.79	1.23	1.33
1	A	367	SER	C-N	-6.50	1.24	1.33
1	C	105	VAL	CA-CB	5.68	1.57	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	104	ASP	CA-C-N	6.73	124.48	120.24
1	C	104	ASP	C-N-CA	6.73	124.48	120.24
1	G	149	ILE	CB-CA-C	-6.58	103.14	112.22
2	F	100	ILE	N-CA-C	6.21	116.99	110.72
2	H	168	ILE	CB-CA-C	-5.09	104.80	110.96
1	C	76	ALA	CA-C-N	5.04	125.03	119.89
1	C	76	ALA	C-N-CA	5.04	125.03	119.89

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	367	SER	Peptide
1	E	292	LYS	Peptide

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	2592	0	2520	26	0
1	C	2659	0	2595	27	0
1	E	2521	0	2455	40	0
1	G	2565	0	2504	15	0
2	B	2428	0	2401	32	0
2	D	2489	0	2515	21	0
2	F	2369	0	2401	25	0
2	H	2416	0	2420	27	0
3	I	303	0	316	11	0
3	J	311	0	318	10	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
5	A	26	0	16	1	0
5	C	26	0	16	0	0
5	E	26	0	16	1	0
5	G	26	0	16	0	0
6	A	96	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	124	0	0	3	0
6	C	168	0	0	1	0
6	D	209	0	0	2	0
6	E	78	0	0	4	0
6	F	194	0	0	3	0
6	G	88	0	0	2	0
6	H	161	0	0	4	0
6	I	24	0	0	4	0
6	J	31	0	0	3	0
All	All	21934	0	20509	209	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 5.

All (209) 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:E:303:LEU:CD1	1:E:309:TRP:HB2	1.73	1.18
1:E:303:LEU:HD11	1:E:309:TRP:HB2	1.30	1.09
1:E:71:ARG:HG2	1:E:71:ARG:HH11	1.19	1.04
3:J:142:LYS:HB2	6:J:2012:HOH:O	1.71	0.91
6:B:2105:HOH:O	3:I:133:PRO:HB3	1.76	0.85
2:F:143:LYS:HE3	6:F:2095:HOH:O	1.78	0.82
1:E:303:LEU:CD1	1:E:309:TRP:CB	2.59	0.80
1:C:367:SER:N	1:C:368:LYS:HA	1.97	0.79
1:E:303:LEU:HD11	1:E:309:TRP:CB	2.12	0.79
3:I:129:VAL:HG21	3:I:146:ILE:CG2	2.13	0.78
1:C:203:ARG:HA	1:C:210:VAL:HG22	1.64	0.78
2:D:143:LYS:HE2	6:D:2110:HOH:O	1.86	0.75
1:A:9:PHE:HZ	1:C:231:ILE:HD11	1.53	0.73
1:E:93:GLU:HB3	6:E:2022:HOH:O	1.88	0.73
1:E:304:GLU:C	1:E:306:LYS:H	1.98	0.71
3:J:144:VAL:HG23	6:J:2012:HOH:O	1.88	0.71
2:F:80:ILE:HG13	2:F:114:ILE:HG23	1.73	0.71
1:C:38:MET:HE1	1:C:239:LEU:HD13	1.74	0.70
1:A:294:ASP:HB3	1:A:297:VAL:HG23	1.73	0.69
2:B:157:LEU:O	2:B:161:ILE:HG13	1.95	0.66
2:B:93:ILE:HA	2:B:97:MET:HG2	1.77	0.66
1:E:71:ARG:HG2	1:E:71:ARG:NH1	1.98	0.66
2:H:82:PHE:HB2	2:H:85:PHE:CZ	2.31	0.66
2:B:59:GLU:HB3	2:B:89:VAL:HG22	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:76:ALA:HB1	1:E:103:ARG:HG2	1.77	0.66
1:E:293:LYS:O	1:E:294:ASP:C	2.38	0.65
1:E:303:LEU:HD12	1:E:309:TRP:HB2	1.72	0.65
3:I:129:VAL:HG21	3:I:146:ILE:HG22	1.78	0.65
2:B:142:LEU:HD23	2:B:166:PRO:HB2	1.78	0.64
2:F:277[B]:ASN:HD21	2:H:277:ASN:HB2	1.64	0.63
2:B:324:PHE:O	3:I:137:LYS:NZ	2.24	0.63
1:E:71:ARG:HH11	1:E:71:ARG:CG	2.06	0.63
2:F:277[B]:ASN:ND2	2:H:277:ASN:HB2	2.14	0.62
1:E:83:ALA:O	1:E:87:ALA:HB3	2.00	0.62
1:C:249:ARG:O	1:C:253:ILE:HG23	1.99	0.61
2:D:214:HIS:HD2	6:D:2147:HOH:O	1.83	0.61
2:B:277[B]:ASN:ND2	2:D:277:ASN:HB2	2.15	0.60
2:F:80:ILE:HG21	2:F:89:VAL:HG11	1.81	0.60
3:I:161:GLU:HB3	6:I:2018:HOH:O	2.01	0.60
1:E:304:GLU:C	1:E:306:LYS:N	2.59	0.59
2:B:264:ARG:NH2	6:B:2096:HOH:O	2.34	0.59
1:E:43:ASP:HA	1:E:46:LEU:HD12	1.85	0.58
1:A:344:ILE:HG23	1:A:355:LEU:HD22	1.85	0.58
3:I:160:LYS:HB3	6:I:2017:HOH:O	2.03	0.58
1:A:142:ILE:HG21	2:D:126:GLU:O	2.05	0.57
1:A:228:ILE:HG21	1:A:258:PRO:HG2	1.86	0.56
1:G:83:ALA:O	1:G:87:ALA:HB3	2.05	0.56
1:G:313:GLU:O	1:G:317:VAL:HG23	2.06	0.56
2:D:45:ALA:O	2:F:3:MET:HE2	2.06	0.56
1:A:249:ARG:O	1:A:253:ILE:HG12	2.06	0.56
2:F:299:PRO:HD3	2:F:307:TRP:CD2	2.40	0.55
3:J:129:VAL:HG21	3:J:146:ILE:CG2	2.37	0.55
3:J:145:ASP:HB3	3:J:148:LEU:HD22	1.88	0.55
2:H:93:ILE:HG21	2:H:168:ILE:HD11	1.89	0.55
1:G:202:ASN:O	1:G:203:ARG:HB2	2.06	0.54
1:C:367:SER:H	1:C:368:LYS:HA	1.72	0.54
2:D:122:VAL:HG23	2:D:124:THR:HG23	1.89	0.54
2:H:282:LEU:HA	3:J:134:SER:HB3	1.89	0.53
2:F:128:HIS:N	6:F:2072:HOH:O	2.42	0.53
2:H:82:PHE:CZ	2:H:173:LEU:HD11	2.43	0.53
2:F:80:ILE:HG13	2:F:114:ILE:CG2	2.38	0.53
2:D:298:TYR:CG	2:D:299:PRO:HD2	2.44	0.53
1:C:23:PHE:HB3	1:C:233:VAL:HG12	1.89	0.53
3:I:129:VAL:HG21	3:I:146:ILE:HG21	1.90	0.53
1:G:66:LEU:HD13	1:G:329:ILE:HD11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:ASP:HB3	1:A:297:VAL:CG2	2.38	0.52
1:C:46:LEU:HB3	1:C:303:LEU:HD21	1.90	0.52
2:H:17:LEU:HD11	2:H:25:ILE:HD12	1.92	0.52
1:A:60:ASP:OD2	1:A:79:ALA:HB2	2.08	0.52
2:H:299:PRO:HD3	2:H:307:TRP:CD2	2.44	0.52
1:A:9:PHE:CZ	1:C:231:ILE:HD11	2.40	0.52
1:E:49:LEU:O	1:E:53:MET:HG3	2.10	0.52
1:A:183:GLN:OE1	2:B:58:ALA:HB1	2.10	0.52
2:F:299:PRO:HA	2:H:102:TYR:OH	2.10	0.52
2:B:284:LEU:HD22	2:B:286:ALA:O	2.10	0.52
2:B:299:PRO:HD3	2:B:307:TRP:CD2	2.45	0.52
1:A:25:ILE:HD12	1:A:240:ALA:HA	1.91	0.52
1:E:304:GLU:O	1:E:307:GLY:N	2.39	0.51
1:A:49:LEU:HD11	1:A:87:ALA:HA	1.93	0.51
1:C:183:GLN:OE1	2:D:58:ALA:HB1	2.10	0.51
2:B:298:TYR:CG	2:B:299:PRO:HD2	2.45	0.51
1:E:23:PHE:HB3	1:E:233:VAL:HG12	1.93	0.51
2:D:93:ILE:HA	2:D:97:MET:HG2	1.92	0.50
1:E:78:THR:HG21	1:E:85:GLN:OE1	2.10	0.50
1:E:294:ASP:C	1:E:294:ASP:OD1	2.54	0.50
2:F:28:GLU:HG2	2:F:57:LEU:HD23	1.94	0.50
2:B:35:GLY:HA3	6:B:2014:HOH:O	2.11	0.50
1:G:67:ASN:ND2	1:G:73:GLY:O	2.44	0.50
2:D:75:ARG:HD3	2:D:161:ILE:O	2.11	0.49
1:G:66:LEU:HD22	1:G:329:ILE:CD1	2.42	0.49
1:A:58:ILE:O	1:A:62:ARG:HG2	2.13	0.49
2:D:174:LYS:O	2:D:178:SER:HB3	2.12	0.49
2:H:82:PHE:CB	2:H:85:PHE:CZ	2.96	0.49
1:C:100:PRO:HG3	1:C:108:ILE:HG21	1.94	0.48
1:E:303:LEU:HD12	1:E:309:TRP:CB	2.33	0.48
2:F:174:LYS:O	2:F:178:SER:HB3	2.13	0.48
1:E:36:GLU:HG2	6:E:2005:HOH:O	2.13	0.48
2:B:142:LEU:CD2	2:B:166:PRO:HB2	2.43	0.48
2:D:117:PRO:HA	2:D:171:GLU:O	2.13	0.48
1:E:25:ILE:HD13	1:E:39:PRO:HD3	1.95	0.48
2:H:298:TYR:CG	2:H:299:PRO:HD2	2.48	0.48
2:F:38:ARG:HB2	2:F:177:ARG:HH22	1.79	0.47
1:A:348:PHE:CZ	2:B:141:GLY:HA2	2.49	0.47
1:A:231:ILE:HD11	1:C:9:PHE:HZ	1.79	0.47
2:B:16:GLU:OE2	2:B:158:ILE:HG23	2.14	0.47
1:E:300:ARG:NH2	1:E:314:GLU:OE2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:27:ASN:HD21	1:G:31:GLU:HB2	1.80	0.47
3:J:141:GLU:HG3	3:J:142:LYS:HG3	1.96	0.47
2:F:116:SER:HB3	2:F:170:LEU:HD23	1.97	0.47
2:B:102:TYR:OH	2:D:299:PRO:HA	2.15	0.47
2:B:81:GLN:O	2:B:117:PRO:HG2	2.15	0.47
1:A:88:SER:HB2	1:A:245:VAL:HG21	1.97	0.46
2:B:122:VAL:HG23	2:B:124:THR:HG23	1.96	0.46
2:B:238:VAL:HB	2:B:268:ILE:HG12	1.96	0.46
1:C:38:MET:HE2	6:C:2014:HOH:O	2.15	0.46
1:C:83:ALA:O	1:C:87:ALA:HB3	2.16	0.46
1:C:367:SER:N	1:C:368:LYS:CA	2.71	0.46
1:E:50:MET:HE1	1:E:300:ARG:HB2	1.98	0.46
2:H:59:GLU:H	2:H:59:GLU:CD	2.24	0.46
1:G:149:ILE:H	1:G:149:ILE:HG13	1.41	0.46
3:J:129:VAL:HG21	3:J:146:ILE:HG22	1.97	0.46
1:E:78:THR:HG21	1:E:85:GLN:CD	2.40	0.46
1:G:298:ARG:HB3	6:G:2060:HOH:O	2.15	0.46
1:A:9:PHE:CD1	1:C:229:PRO:HB3	2.52	0.46
1:A:368:LYS:N	6:A:2095:HOH:O	2.48	0.45
2:F:74:PHE:O	2:F:76:PRO:HD3	2.15	0.45
2:H:82:PHE:HZ	2:H:173:LEU:HD11	1.81	0.45
2:H:185:GLU:HG3	6:H:2078:HOH:O	2.15	0.45
2:B:20:ASP:HB3	2:B:23:VAL:HG23	1.97	0.45
6:E:2037:HOH:O	2:F:96:GLN:HG3	2.16	0.45
1:E:60:ASP:OD2	1:E:79:ALA:HB2	2.16	0.45
1:E:144:ILE:HD12	5:E:1370:TPP:C2'	2.47	0.45
2:F:59:GLU:CD	2:F:59:GLU:H	2.25	0.45
1:C:51:ARG:HG2	1:C:317:VAL:HG21	1.98	0.45
2:B:128:HIS:HE1	1:C:142:ILE:HD12	1.82	0.44
1:E:42:SER:H	1:E:45[B]:GLN:HE21	1.64	0.44
2:B:36:VAL:HG13	2:B:37:PHE:CD2	2.53	0.44
1:A:187:PHE:HE1	2:B:58:ALA:HB3	1.81	0.44
2:B:128:HIS:HE1	1:C:142:ILE:CD1	2.31	0.44
2:H:38:ARG:HG2	6:H:2015:HOH:O	2.16	0.44
2:H:114:ILE:HD13	2:H:168:ILE:HD13	1.99	0.44
2:H:286:ALA:N	3:J:132:MET:HE2	2.33	0.44
1:C:203:ARG:NH2	1:C:264:LEU:HD22	2.33	0.43
2:D:299:PRO:HD3	2:D:307:TRP:CD2	2.53	0.43
1:G:26:LEU:HG	1:G:239:LEU:HD12	2.00	0.43
1:A:100:PRO:HG2	1:A:140:PRO:HA	2.00	0.43
1:A:224:VAL:HG13	1:C:16:VAL:HG11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1370:TPP:HM42	2:D:57:LEU:HD11	1.99	0.43
1:C:38:MET:HE3	1:C:39:PRO:HD2	2.00	0.43
2:D:120:GLY:HA3	2:D:298:TYR:OH	2.18	0.43
1:E:293:LYS:O	1:E:294:ASP:O	2.35	0.43
1:E:326:LYS:O	1:E:329:ILE:HG22	2.18	0.43
2:H:205:THR:O	2:H:256:ALA:HA	2.19	0.43
2:B:59:GLU:HB3	2:B:89:VAL:CG2	2.47	0.43
2:F:264:ARG:NH1	6:F:2156:HOH:O	2.33	0.43
2:H:180:ARG:HD2	6:H:2004:HOH:O	2.19	0.43
2:D:125:PRO:HG2	2:D:128:HIS:CE1	2.53	0.43
1:E:38:MET:HE3	1:E:239:LEU:HD13	2.01	0.43
1:C:310:SER:HG	1:C:313:GLU:HG3	1.84	0.43
3:I:129:VAL:HG13	6:I:2008:HOH:O	2.19	0.42
1:A:83:ALA:O	1:A:87:ALA:HB3	2.19	0.42
2:B:76:PRO:HD2	2:B:112:ILE:HG12	2.01	0.42
2:H:93:ILE:CG2	2:H:168:ILE:HD11	2.49	0.42
1:C:26:LEU:HG	1:C:239:LEU:HD12	2.00	0.42
3:J:144:VAL:HG12	6:J:2020:HOH:O	2.19	0.42
2:B:284:LEU:O	3:I:136:ARG:HD2	2.19	0.42
1:C:203:ARG:O	1:C:209:PRO:HA	2.19	0.42
2:D:286:ALA:HB1	2:D:287:PRO:HD2	2.00	0.42
1:G:123:ARG:NH2	1:G:333:ASP:OD1	2.46	0.42
6:G:2039:HOH:O	2:H:96:GLN:HG3	2.19	0.42
2:H:14:ARG:HD3	6:H:2018:HOH:O	2.19	0.42
1:E:304:GLU:O	1:E:306:LYS:N	2.53	0.42
2:F:282:LEU:HD13	3:J:130:ILE:HB	2.02	0.42
2:H:162:ARG:HH22	2:H:187:GLU:CD	2.27	0.42
1:A:298:ARG:HB3	6:A:2071:HOH:O	2.19	0.42
2:F:93:ILE:HA	2:F:97:MET:CG	2.48	0.42
2:H:80:ILE:O	2:H:80:ILE:HG22	2.18	0.42
2:D:20:ASP:HB3	2:D:23:VAL:HG23	2.02	0.42
1:G:66:LEU:HD22	1:G:329:ILE:HD11	2.02	0.42
2:B:117:PRO:HA	2:B:171:GLU:O	2.19	0.42
1:A:105:VAL:N	1:A:106:PRO:CD	2.82	0.41
2:F:17:LEU:HD11	2:F:25:ILE:HD12	2.02	0.41
2:B:299:PRO:HD3	2:B:307:TRP:CE3	2.54	0.41
2:D:298:TYR:CD2	2:D:299:PRO:HD2	2.56	0.41
1:A:237:ASP:OD1	1:A:237:ASP:C	2.63	0.41
2:F:97:MET:SD	2:F:114:ILE:HD11	2.60	0.41
2:F:142:LEU:HD23	2:F:166:PRO:HB2	2.01	0.41
1:E:294:ASP:HA	1:E:295:PRO:HD2	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:100:ILE:HD13	2:B:110:MET:HE3	2.02	0.41
1:C:203:ARG:HH21	1:C:264:LEU:HD22	1.85	0.41
2:D:97:MET:HA	2:D:100:ILE:HD12	2.02	0.41
1:E:235:GLY:N	1:E:264:LEU:O	2.49	0.41
1:E:338:GLN:HE21	2:H:302:GLN:HB3	1.85	0.41
1:G:335:THR:HA	1:G:336:PRO:HD3	1.97	0.41
2:F:300:PHE:O	2:F:301:ALA:C	2.64	0.41
1:E:76:ALA:CB	1:E:103:ARG:HG2	2.47	0.40
1:E:83:ALA:HB3	1:E:235:GLY:O	2.21	0.40
3:I:132:MET:HE3	3:I:133:PRO:HD2	2.03	0.40
1:A:50:MET:O	1:A:54:VAL:HG23	2.21	0.40
2:B:38:ARG:HG3	2:B:177:ARG:NH2	2.36	0.40
1:E:142:ILE:HG13	6:E:2036:HOH:O	2.22	0.40
2:F:92:SER:O	2:F:97:MET:HG2	2.22	0.40
2:H:142:LEU:HD22	2:H:168:ILE:HG12	2.03	0.40
1:E:7:PHE:O	1:E:9:PHE:N	2.54	0.40
1:G:346:ILE:HD12	2:H:102:TYR:O	2.21	0.40
3:I:161:GLU:HB2	6:I:2021:HOH:O	2.21	0.40
2:B:97:MET:SD	2:B:114:ILE:HD11	2.61	0.40
1:C:51:ARG:HG2	1:C:317:VAL:CG2	2.52	0.40
1:G:199:VAL:HG23	1:G:260:LEU:HD11	2.02	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	337/368 (92%)	327 (97%)	9 (3%)	1 (0%)	36	46
1	C	339/368 (92%)	329 (97%)	9 (3%)	1 (0%)	36	46
1	E	323/368 (88%)	304 (94%)	16 (5%)	3 (1%)	14	17
1	G	326/368 (89%)	316 (97%)	9 (3%)	1 (0%)	36	46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	323/324 (100%)	309 (96%)	13 (4%)	1 (0%)	36	46
2	D	322/324 (99%)	311 (97%)	11 (3%)	0	100	100
2	F	305/324 (94%)	298 (98%)	6 (2%)	1 (0%)	36	46
2	H	313/324 (97%)	303 (97%)	8 (3%)	2 (1%)	21	27
3	I	38/49 (78%)	38 (100%)	0	0	100	100
3	J	41/49 (84%)	40 (98%)	1 (2%)	0	100	100
All	All	2667/2866 (93%)	2575 (97%)	82 (3%)	10 (0%)	30	38

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	ILE
1	C	206	ILE
1	E	305	ALA
1	G	203	ARG
2	H	81	GLN
2	F	301	ALA
1	E	297	VAL
2	B	184	PRO
2	H	86	VAL
1	E	294	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	255/304 (84%)	251 (98%)	4 (2%)	55	73
1	C	267/304 (88%)	259 (97%)	8 (3%)	36	53
1	E	250/304 (82%)	240 (96%)	10 (4%)	28	42
1	G	257/304 (84%)	255 (99%)	2 (1%)	73	86
2	B	247/263 (94%)	238 (96%)	9 (4%)	31	47
2	D	263/263 (100%)	258 (98%)	5 (2%)	50	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	249/263 (95%)	240 (96%)	9 (4%)	31	47
2	H	251/263 (95%)	248 (99%)	3 (1%)	63	79
3	I	31/36 (86%)	27 (87%)	4 (13%)	4	4
3	J	30/36 (83%)	26 (87%)	4 (13%)	4	4
All	All	2100/2340 (90%)	2042 (97%)	58 (3%)	39	56

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	142	ILE
1	A	265	CYS
1	A	293	LYS
1	A	329	ILE
2	B	38	ARG
2	B	89	VAL
2	B	100	ILE
2	B	131	SER
2	B	176	TYR
2	B	184	PRO
2	B	187	GLU
2	B	277[A]	ASN
2	B	277[B]	ASN
1	C	13	LEU
1	C	142	ILE
1	C	203	ARG
1	C	210	VAL
1	C	211	GLU
1	C	253	ILE
1	C	254	ASN
1	C	265	CYS
2	D	100	ILE
2	D	127	LEU
2	D	187	GLU
2	D	277	ASN
2	D	320	LYS
1	E	38	MET
1	E	67	ASN
1	E	68	ARG
1	E	71	ARG
1	E	78	THR
1	E	137	VAL

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Mol	Chain	Res	Type
1	E	202	ASN
1	E	303	LEU
1	E	329	ILE
1	E	368	LYS
2	F	32	VAL
2	F	80	ILE
2	F	89	VAL
2	F	97	MET
2	F	187	GLU
2	F	277[A]	ASN
2	F	277[B]	ASN
2	F	304	GLU
2	F	320	LYS
1	G	149	ILE
1	G	329	ILE
2	H	131	SER
2	H	142	LEU
2	H	277	ASN
3	I	129	VAL
3	I	146	ILE
3	I	157	ARG
3	I	160	LYS
3	J	129	VAL
3	J	146	ILE
3	J	148	LEU
3	J	160	LYS

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

Mol	Chain	Res	Type
1	A	24	GLN
1	A	69	GLN
2	B	128	HIS
1	C	67	ASN
1	C	200	GLN
2	D	123	HIS
2	D	271	ASN
1	E	19	GLN
1	E	24	GLN
1	E	81	GLN
1	E	183	GLN
1	E	316	ASN

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Mol	Chain	Res	Type
1	E	320	GLN
2	F	271	ASN
2	F	323	ASN
1	G	67	ASN
1	G	81	GLN
1	G	178	GLN
1	G	186	ASN
2	H	271	ASN
2	H	302	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	TPP	C	1370	4	26,27,27	2.18	4 (15%)	38,40,40	1.63	11 (28%)
5	TPP	E	1370	4	26,27,27	2.34	6 (23%)	38,40,40	1.59	9 (23%)
5	TPP	G	1370	4	26,27,27	2.20	4 (15%)	38,40,40	1.46	7 (18%)
5	TPP	A	1370	4	26,27,27	2.28	5 (19%)	38,40,40	1.52	10 (26%)

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	C	1370	4	-	1/17/17/17	0/2/2/2
5	TPP	E	1370	4	-	4/17/17/17	0/2/2/2
5	TPP	G	1370	4	-	8/17/17/17	0/2/2/2
5	TPP	A	1370	4	-	1/17/17/17	0/2/2/2

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	1370	TPP	PA-O3A	-7.45	1.51	1.59
5	C	1370	TPP	C5-C4	6.81	1.49	1.35
5	A	1370	TPP	C5-C4	6.71	1.49	1.35
5	E	1370	TPP	C5-C4	6.67	1.49	1.35
5	G	1370	TPP	C5-C4	6.58	1.49	1.35
5	G	1370	TPP	PA-O3A	-6.39	1.52	1.59
5	A	1370	TPP	PA-O3A	-6.04	1.53	1.59
5	C	1370	TPP	PA-O3A	-5.80	1.53	1.59
5	A	1370	TPP	C5'-C4'	4.75	1.50	1.42
5	C	1370	TPP	C5'-C4'	4.22	1.49	1.42
5	E	1370	TPP	C5'-C4'	4.10	1.49	1.42
5	G	1370	TPP	C5'-C4'	4.09	1.49	1.42
5	A	1370	TPP	C4'-N4'	2.83	1.41	1.34
5	E	1370	TPP	C4'-N4'	2.55	1.40	1.34
5	G	1370	TPP	C4'-N4'	2.50	1.40	1.34
5	C	1370	TPP	C4'-N4'	2.47	1.40	1.34
5	E	1370	TPP	C4-N3	-2.07	1.35	1.39
5	E	1370	TPP	PA-O7	-2.02	1.51	1.59
5	A	1370	TPP	C4-N3	-2.01	1.35	1.39

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	1370	TPP	C6'-N1'-C2'	3.84	122.38	116.07
5	A	1370	TPP	C6'-N1'-C2'	3.60	121.98	116.07
5	G	1370	TPP	C6'-N1'-C2'	3.59	121.97	116.07
5	C	1370	TPP	C6'-N1'-C2'	3.59	121.96	116.07
5	C	1370	TPP	C6-C5-S1	-3.10	113.64	120.90
5	C	1370	TPP	CM4-C4-N3	2.87	127.19	120.57
5	E	1370	TPP	S1-C2-N3	2.81	115.85	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	1370	TPP	C5'-C6'-N1'	-2.71	119.43	123.83
5	E	1370	TPP	C6-C5-S1	-2.70	114.58	120.90
5	G	1370	TPP	C5'-C6'-N1'	-2.66	119.51	123.83
5	A	1370	TPP	C6-C5-S1	-2.63	114.75	120.90
5	A	1370	TPP	O2B-PB-O3A	2.61	113.40	104.64
5	G	1370	TPP	CM4-C4-N3	2.61	126.59	120.57
5	G	1370	TPP	S1-C2-N3	2.57	115.55	112.30
5	C	1370	TPP	C6-C5-C4	2.55	134.58	128.17
5	A	1370	TPP	N1'-C2'-N3'	-2.48	121.40	125.53
5	E	1370	TPP	C2-S1-C5	-2.48	89.57	91.22
5	C	1370	TPP	N1'-C2'-N3'	-2.48	121.41	125.53
5	E	1370	TPP	N1'-C2'-N3'	-2.45	121.46	125.53
5	E	1370	TPP	CM4-C4-N3	2.44	126.19	120.57
5	E	1370	TPP	N4'-C4'-N3'	2.37	120.22	117.03
5	C	1370	TPP	S1-C2-N3	2.35	115.27	112.30
5	A	1370	TPP	S1-C2-N3	2.27	115.17	112.30
5	A	1370	TPP	C5'-C6'-N1'	-2.26	120.15	123.83
5	G	1370	TPP	N4'-C4'-N3'	2.22	120.02	117.03
5	A	1370	TPP	CM4-C4-N3	2.22	125.68	120.57
5	C	1370	TPP	C5'-C6'-N1'	-2.20	120.25	123.83
5	C	1370	TPP	CM2-C2'-N1'	2.18	119.52	117.20
5	C	1370	TPP	C7-C6-C5	-2.14	106.02	112.73
5	A	1370	TPP	C5'-C4'-N3'	-2.13	118.18	121.31
5	G	1370	TPP	N1'-C2'-N3'	-2.12	122.00	125.53
5	G	1370	TPP	C5'-C4'-N3'	-2.12	118.20	121.31
5	A	1370	TPP	C2'-N3'-C4'	2.07	121.28	118.10
5	C	1370	TPP	O2B-PB-O3A	2.06	111.56	104.64
5	C	1370	TPP	C2'-N3'-C4'	2.04	121.23	118.10
5	A	1370	TPP	C6-C5-C4	2.01	133.22	128.17
5	E	1370	TPP	C6-C5-C4	2.01	133.21	128.17

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	1370	TPP	PB-O3A-PA-O7
5	E	1370	TPP	C5'-C7'-N3-C2
5	E	1370	TPP	C5-C6-C7-O7
5	E	1370	TPP	PB-O3A-PA-O7
5	G	1370	TPP	C5-C6-C7-O7
5	G	1370	TPP	C7-O7-PA-O1A
5	G	1370	TPP	C7-O7-PA-O2A

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Mol	Chain	Res	Type	Atoms
5	E	1370	TPP	C5'-C7'-N3-C4
5	A	1370	TPP	PB-O3A-PA-O7
5	G	1370	TPP	C7-O7-PA-O3A
5	G	1370	TPP	S1-C5-C6-C7
5	G	1370	TPP	C4-C5-C6-C7
5	G	1370	TPP	PA-O3A-PB-O3B
5	G	1370	TPP	PA-O3A-PB-O1B

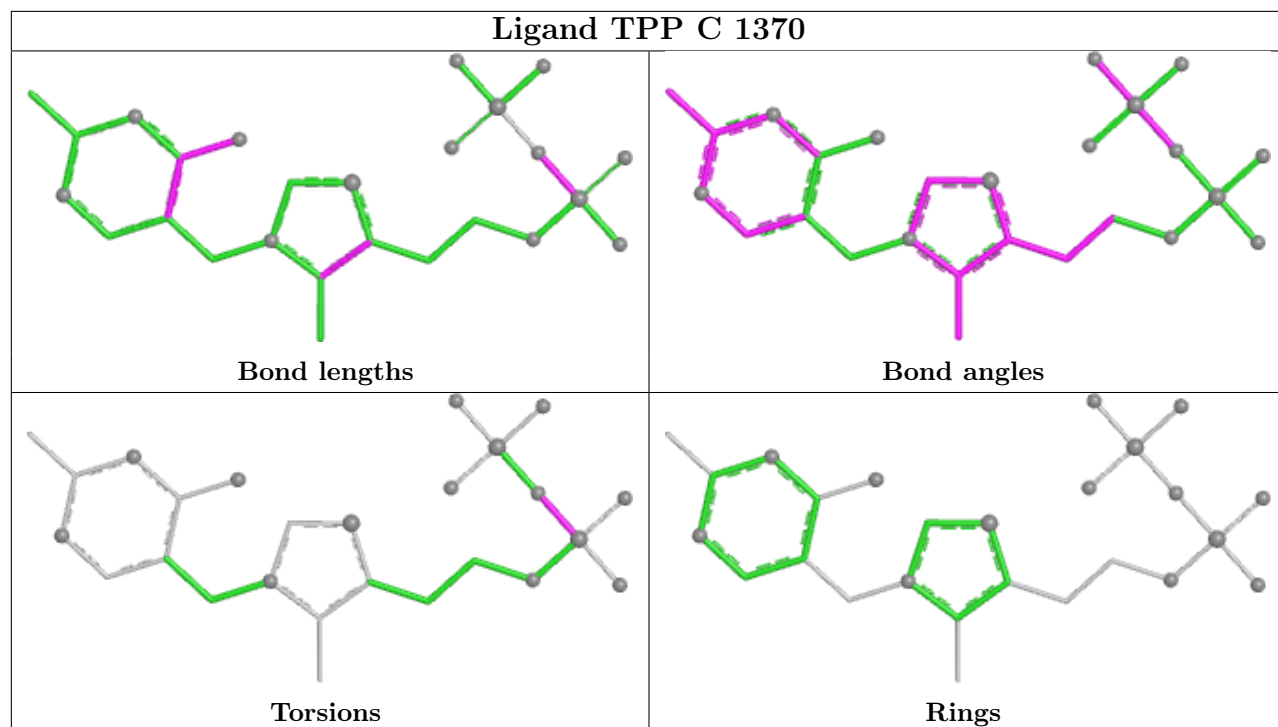
There are no ring outliers.

2 monomers are involved in 2 short contacts:

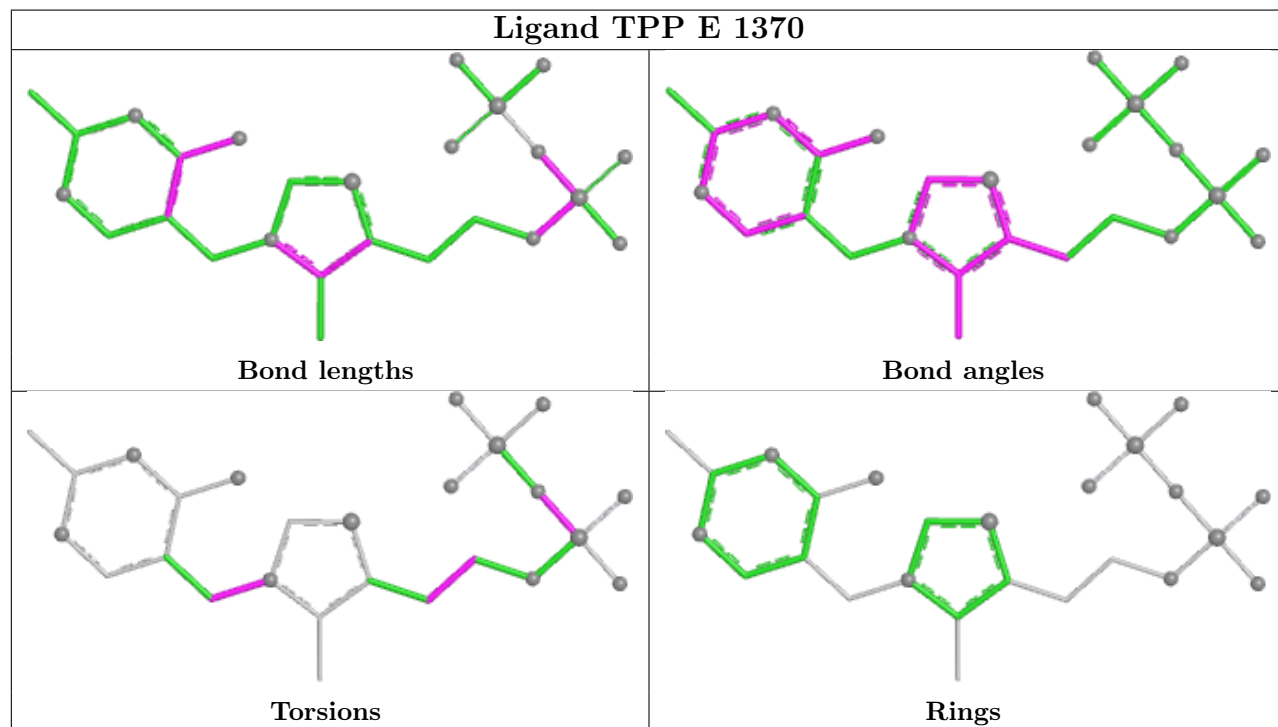
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	1370	TPP	1	0
5	A	1370	TPP	1	0

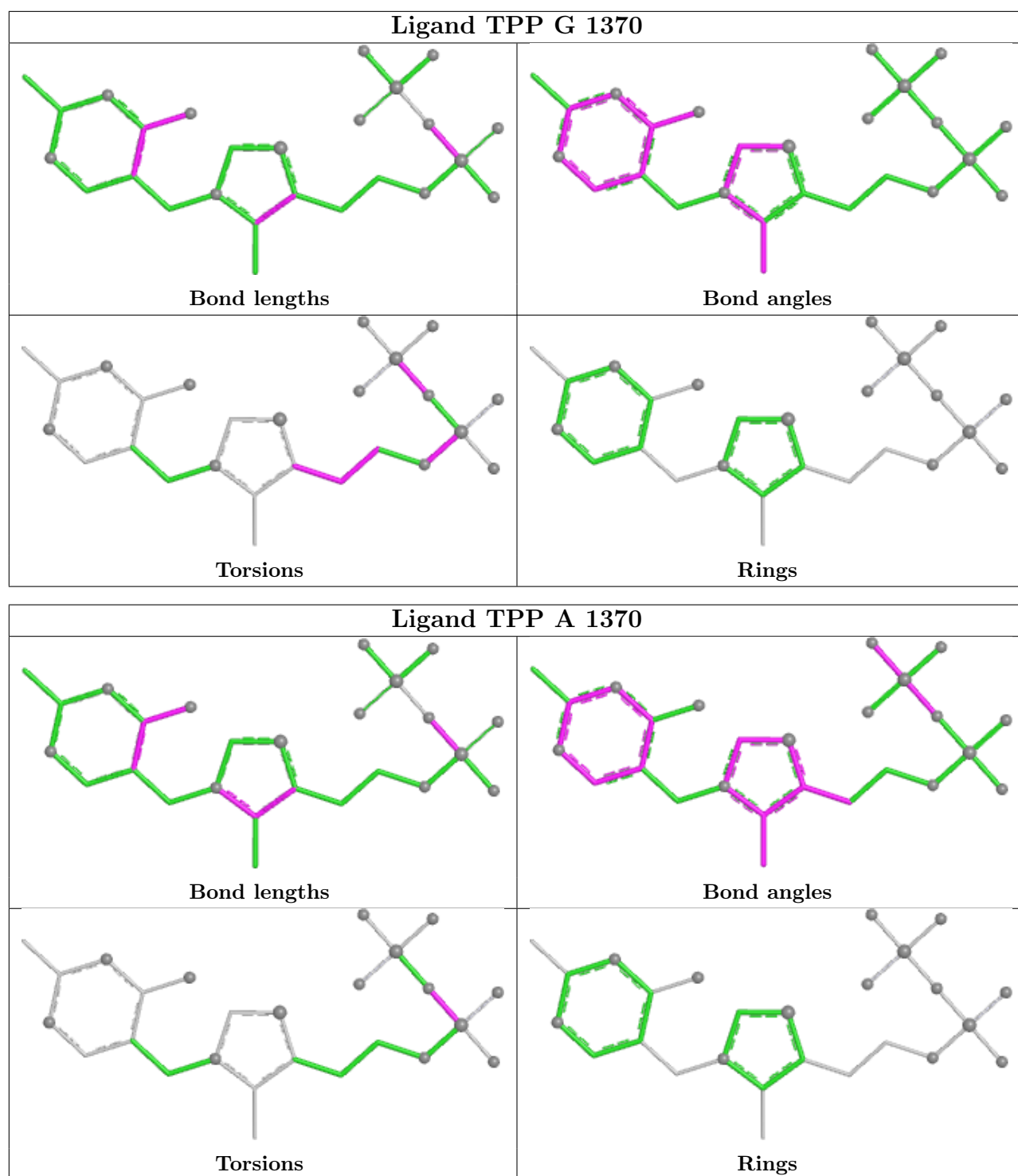
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand TPP C 1370



Ligand TPP E 1370





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 ⓘ

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	341/368 (92%)	0.06	5 (1%) 72 73	24, 42, 62, 79	2 (0%)
1	C	343/368 (93%)	-0.14	6 (1%) 69 70	22, 38, 60, 88	2 (0%)
1	E	328/368 (89%)	0.32	11 (3%) 48 50	20, 47, 68, 85	4 (1%)
1	G	331/368 (89%)	0.06	7 (2%) 63 65	25, 45, 68, 88	2 (0%)
2	B	324/324 (100%)	-0.04	4 (1%) 76 77	20, 37, 53, 67	1 (0%)
2	D	324/324 (100%)	-0.34	2 (0%) 85 86	20, 30, 45, 61	0
2	F	310/324 (95%)	-0.40	5 (1%) 70 72	17, 30, 47, 67	1 (0%)
2	H	317/324 (97%)	-0.24	10 (3%) 50 52	18, 33, 54, 93	0
3	I	40/49 (81%)	1.50	12 (30%) 1 1	50, 66, 76, 82	0
3	J	43/49 (87%)	1.71	16 (37%) 1 1	46, 66, 80, 87	0
All	All	2701/2866 (94%)	-0.03	78 (2%) 53 56	17, 38, 65, 93	12 (0%)

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	51	ARG	5.4
2	H	57	LEU	4.5
1	A	254	ASN	4.3
3	J	160	LYS	4.2
1	C	266	PHE	3.9
2	H	83	PHE	3.9
3	I	163	ILE	3.9
1	E	233	VAL	3.9
1	G	205	ALA	3.8
1	E	204	PHE	3.6
1	E	205	ALA	3.6
2	H	82	PHE	3.3
2	H	86	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	G	233	VAL	3.3
3	J	135	VAL	3.3
2	H	84	GLY	3.2
2	H	121	GLY	3.1
2	B	178	SER	3.1
1	C	5	PHE	3.1
2	H	81	GLN	3.0
1	C	290	TRP	3.0
1	G	204	PHE	2.9
2	H	85	PHE	2.9
3	J	144	VAL	2.9
3	J	139	ALA	2.8
3	I	161	GLU	2.8
1	C	268	TYR	2.8
1	E	303	LEU	2.7
3	J	166	PHE	2.7
3	I	146	ILE	2.6
1	A	266	PHE	2.6
3	J	146	ILE	2.6
1	G	213	GLN	2.6
2	F	57	LEU	2.6
3	I	166	PHE	2.6
2	H	120	GLY	2.6
3	J	138	TYR	2.6
2	F	87	TYR	2.6
1	C	233	VAL	2.5
1	E	292	LYS	2.5
1	A	233	VAL	2.5
3	J	149	VAL	2.5
3	J	148	LEU	2.5
1	G	74	PHE	2.4
3	I	130	ILE	2.4
3	J	132	MET	2.4
1	E	5	PHE	2.4
3	I	144	VAL	2.4
3	J	130	ILE	2.4
3	I	159	LEU	2.4
1	E	254	ASN	2.4
3	I	160	LYS	2.3
2	F	91	ASP	2.3
2	B	186	GLY	2.3
3	I	138	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
3	I	139	ALA	2.3
1	G	202	ASN	2.2
2	B	126	GLU	2.2
2	D	127	LEU	2.2
1	A	268	TYR	2.2
2	B	127	LEU	2.2
3	I	149	VAL	2.2
1	E	301	LYS	2.2
1	G	292	LYS	2.2
3	J	170	GLY	2.2
1	E	78	THR	2.2
1	A	292	LYS	2.1
2	F	59	GLU	2.1
3	J	158	VAL	2.1
2	F	300	PHE	2.1
3	I	165	ALA	2.1
3	J	168	ALA	2.1
1	C	205	ALA	2.0
2	D	126	GLU	2.0
2	H	91	ASP	2.0
1	E	74	PHE	2.0
3	J	150	GLN	2.0
3	J	163	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

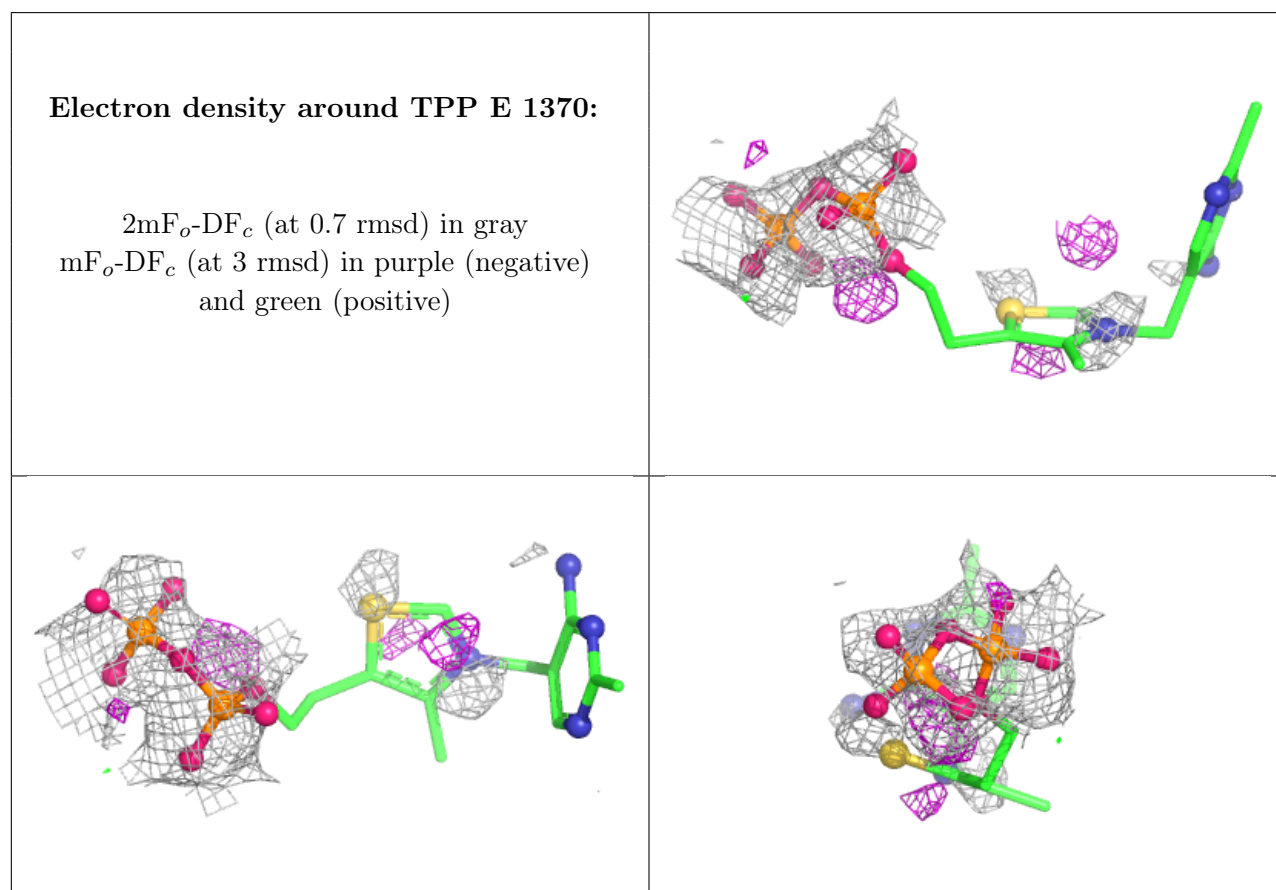
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

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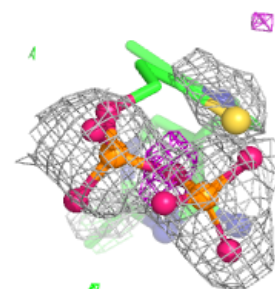
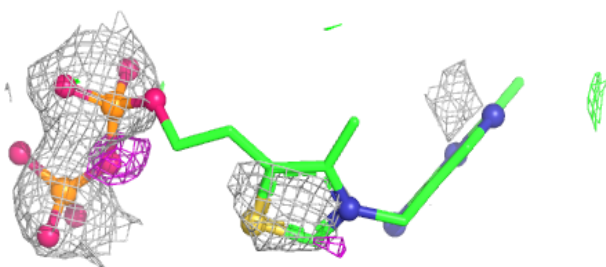
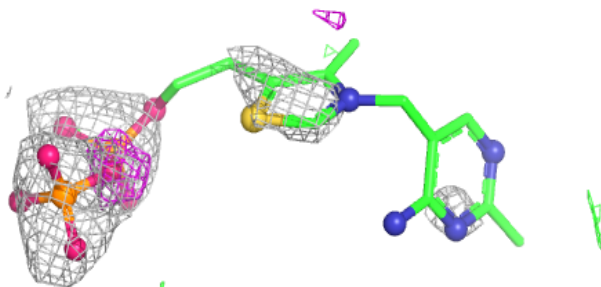
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	TPP	E	1370	26/26	0.63	0.20	172,176,177,178	0
5	TPP	G	1370	26/26	0.63	0.20	143,149,155,155	0
4	MG	E	1368	1/1	0.83	0.17	78,78,78,78	0
4	MG	G	1368	1/1	0.88	0.07	66,66,66,66	0
5	TPP	C	1370	26/26	0.92	0.10	34,47,51,53	0
4	MG	C	1368	1/1	0.94	0.06	28,28,28,28	0
5	TPP	A	1370	26/26	0.95	0.08	29,37,40,42	0
4	MG	A	1368	1/1	0.97	0.04	28,28,28,28	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.

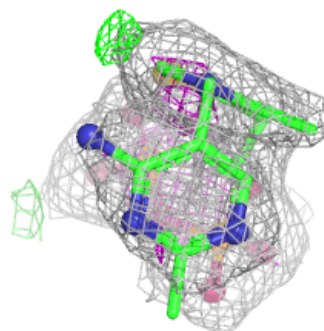
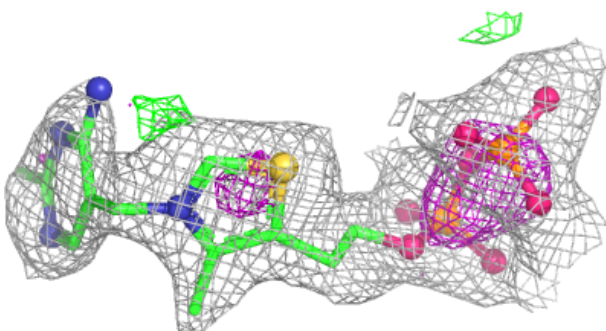
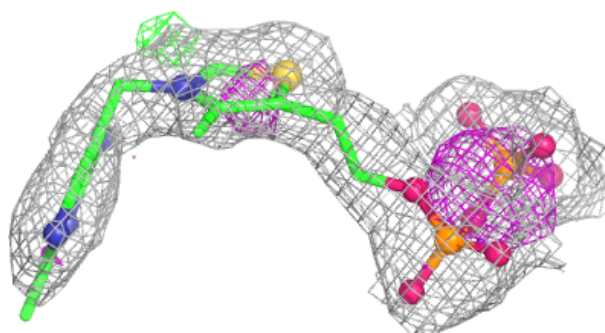


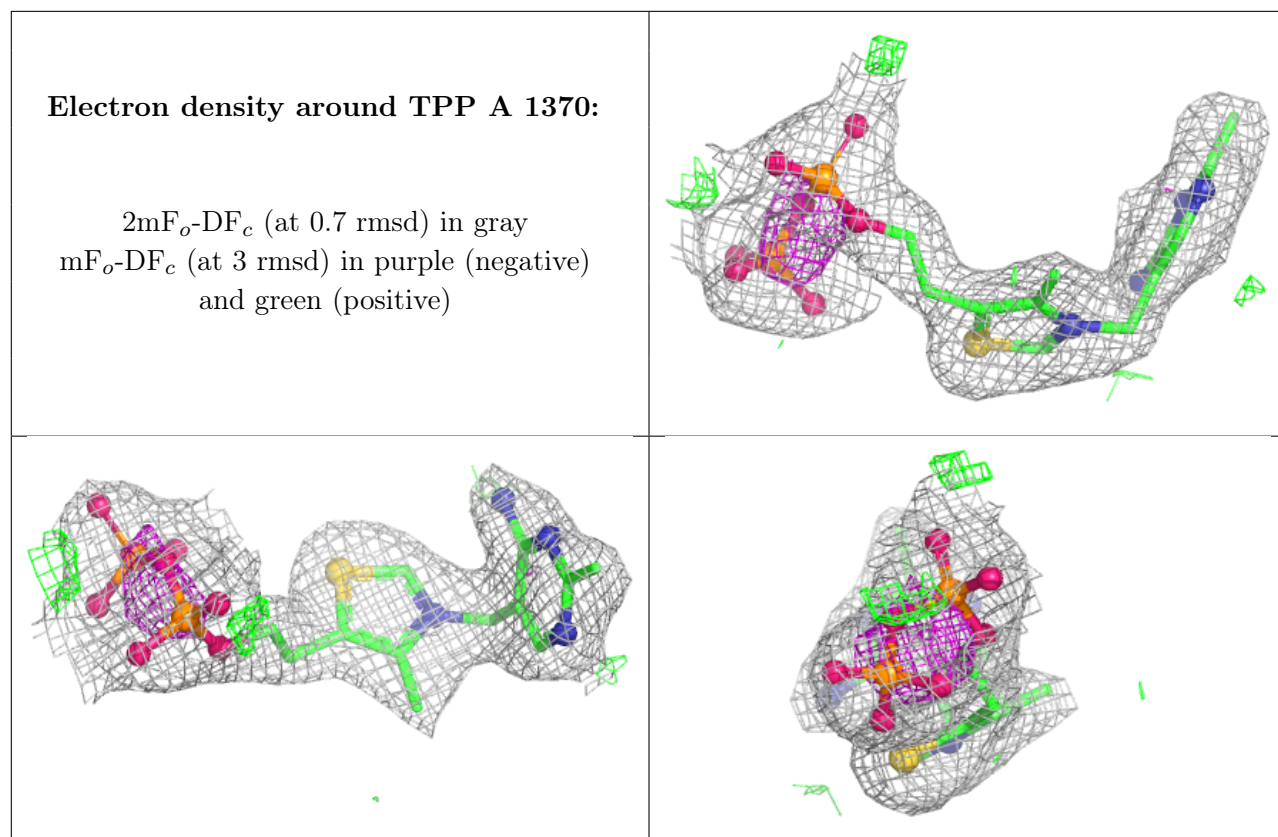
Electron density around TPP G 1370:

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

**Electron density around TPP C 1370:**

$2mF_o-DF_c$ (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.