



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 15, 2020 – 08:34 PM JST

PDB ID : 6L0Q
Title : Crystal Structure of the O-Phosphoserine Sulfhydrylase from Aeropyrum
pernix Complexed with O-Phosphoserine
Authors : Nakabayashi, M.; Takeda, E.; Ishikawa, K.; Nakamura, T.
Deposited on : 2019-09-26
Resolution : 1.58 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

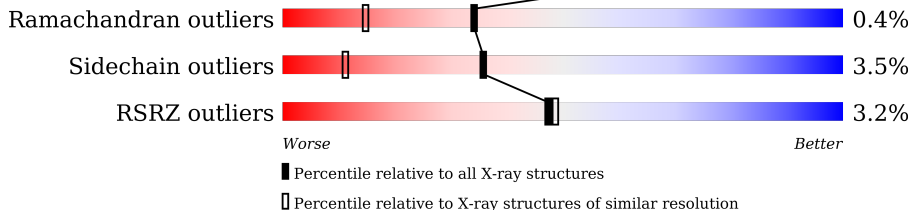
MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.4
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.4

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X-RAY DIFFRACTION

A.

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}	130704	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

fit to the electron density. The red, orange, yellow and green segments on 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	389	<div><div></div><div>4%</div><div>87%</div><div>11%</div><div></div></div>	..	
1	B	389	<div><div></div><div>3%</div><div>86%</div><div>11%</div><div></div></div>	..	
1	C	389	<div><div></div><div>4%</div><div>85%</div><div>13%</div><div></div></div>	..	
1	D	389	<div><div></div><div>2%</div><div>86%</div><div>12%</div><div></div></div>	..	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12171 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.

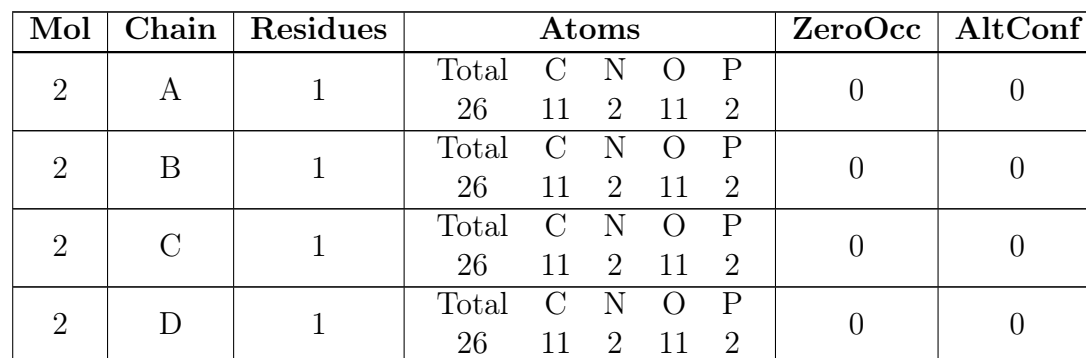
- Molecule 1 is a protein called Protein CysO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	4	0
			2936	1867	514	546	9			
1	B	382	Total	C	N	O	S	0	9	0
			2960	1880	517	554	9			
1	C	382	Total	C	N	O	S	0	5	0
			2938	1865	512	552	9			
1	D	382	Total	C	N	O	S	0	0	0
			2912	1848	509	546	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	ALA	LYS	engineered mutation	UNP Q9YBL2
A	225	TYR	PHE	engineered mutation	UNP Q9YBL2
B	127	ALA	LYS	engineered mutation	UNP Q9YBL2
B	225	TYR	PHE	engineered mutation	UNP Q9YBL2
C	127	ALA	LYS	engineered mutation	UNP Q9YBL2
C	225	TYR	PHE	engineered mutation	UNP Q9YBL2
D	127	ALA	LYS	engineered mutation	UNP Q9YBL2
D	225	TYR	PHE	engineered mutation	UNP Q9YBL2

- Molecule 2 is (2S)-2-[(E)-[2-methyl-3-oxidanyl-5-(phosphonooxymethyl)pyridin-4-yl]methylideneamino]-3-phosphonooxy-propanoic acid (three-letter code: E1U) (formula: C₁₁H₁₆N₂O₁₁P₂) (labeled as "Ligand of Interest" by author).



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- Chemical structure of 2-methylpentan-2-ol (MPD) is shown. The structure is a 5-carbon chain with a methyl group (C5) and a hydroxyl group (OH, O4) attached to the second carbon (C2). The carbon atoms are labeled C1, C2, C3, C4(S), and C5. The hydroxyl group is labeled O2 and OH. The methyl group is labeled CM. The structure is drawn with a central carbon (C2) bonded to a methyl group (C5), a hydroxyl group (OH, O4), and a hydrogen atom (H, O2). The chain continues from C2 to C1, C3, and C4(S).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			8	6	2		
3	C	1	Total	C	O	0	0
			8	6	2		

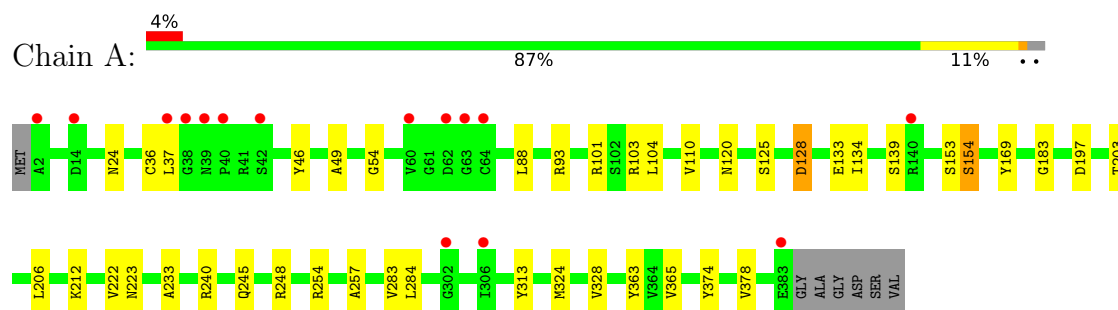
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	61	Total	O	0	0
			61	61		
4	B	87	Total	O	0	0
			87	87		
4	C	75	Total	O	0	0
			75	75		
4	D	82	Total	O	0	0
			82	82		

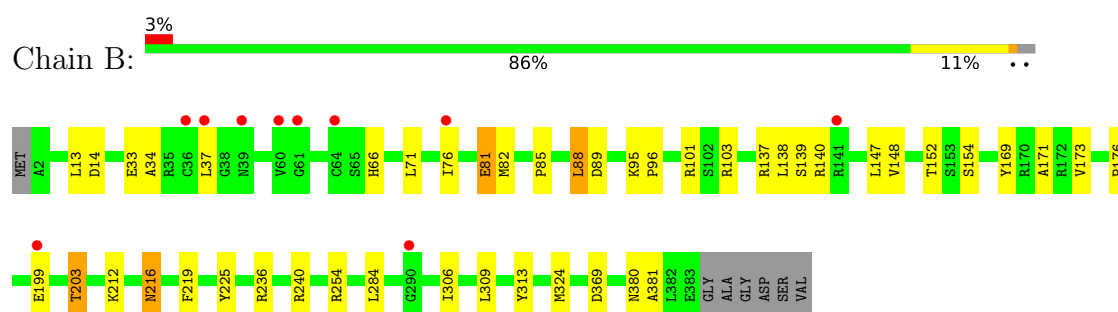
3 Residue-property plots [i](#)

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.

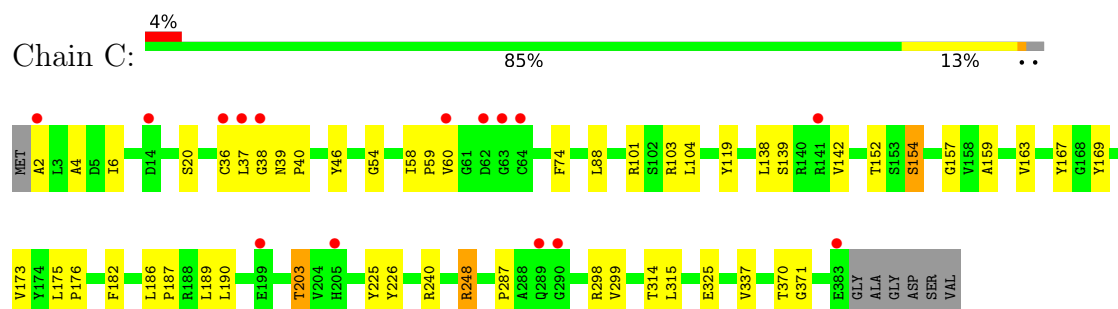
• Molecule 1: Protein CysO



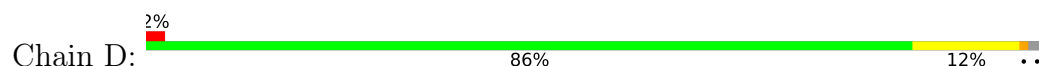
• Molecule 1: Protein CysO

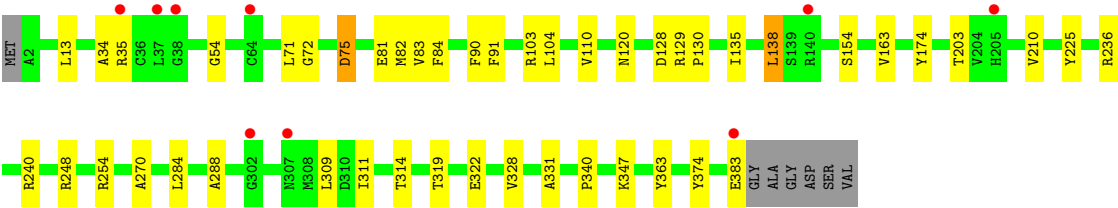


• Molecule 1: Protein CysO



• Molecule 1: Protein CysO





4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	75.20Å 75.20Å 275.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.20 – 1.58 68.82 – 1.58	Depositor EDS
% Data completeness (in resolution range)	100.0 (75.20-1.58) 99.9 (68.82-1.58)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.88 (at 1.58Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.167 , 0.174 0.175 , 0.180	Depositor DCC
R_{free} test set	10205 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.166 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12171	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MPD, E1U

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	0.57	0/3008	0.84	3/4085 (0.1%)
1	B	0.62	0/3047	0.91	4/4137 (0.1%)
1	C	0.55	0/3012	0.83	0/4089
1	D	0.62	0/2971	0.90	3/4035 (0.1%)
All	All	0.59	0/12038	0.87	10/16346 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	254	ARG	NE-CZ-NH2	-6.74	116.93	120.30
1	D	254	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	D	138	LEU	CA-CB-CG	6.16	129.46	115.30
1	B	369	ASP	CB-CG-OD1	-5.72	113.15	118.30
1	A	93	ARG	NE-CZ-NH2	5.65	123.12	120.30
1	B	254	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	A	128	ASP	CB-CG-OD2	-5.25	113.57	118.30
1	B	138	LEU	CA-CB-CG	5.16	127.17	115.30
1	B	101	ARG	NE-CZ-NH2	5.10	122.85	120.30
1	A	103	ARG	NE-CZ-NH1	5.09	122.84	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2936	0	2964	22	0
1	B	2960	0	2991	24	0
1	C	2938	0	2959	27	0
1	D	2912	0	2924	19	0
2	A	26	0	0	0	0
2	B	26	0	0	1	0
2	C	26	0	0	1	0
2	D	26	0	0	1	0
3	B	8	0	14	4	0
3	C	8	0	14	2	0
4	A	61	0	0	0	0
4	B	87	0	0	0	0
4	C	75	0	0	0	0
4	D	82	0	0	0	0
All	All	12171	0	11866	89	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 4.

All (89) 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:37[A]:LEU:HD21	1:A:46:TYR:CD2	2.29	0.68
1:C:159:ALA:O	1:C:163:VAL:HG23	1.96	0.65
1:D:81:GLU:HG2	1:D:83:VAL:HG23	1.82	0.61
1:C:4:ALA:HB2	1:C:60:VAL:HG21	1.83	0.61
1:D:54:GLY:HA3	1:D:240:ARG:HD3	1.83	0.60
3:C:402:MPD:H52	1:D:82:MET:HE1	1.84	0.58
1:B:139:SER:HA	1:B:169:TYR:OH	2.03	0.58
1:A:37[B]:LEU:HD11	1:A:46:TYR:CD2	2.39	0.58
1:C:74:PHE:O	1:C:248:ARG:NH2	2.37	0.57
1:C:88:LEU:HD11	1:C:167:TYR:CE1	2.39	0.57
1:C:54:GLY:HA3	1:C:240:ARG:HD3	1.87	0.56
1:D:225:TYR:OH	2:D:401:E1U:O7P	2.24	0.55
1:B:225:TYR:OH	2:B:401:E1U:O6P	2.23	0.54
1:B:85:PRO:HD2	1:B:89:ASP:OD2	2.08	0.54
1:A:154:SER:OG	1:A:183:GLY:HA3	2.08	0.53
1:D:174:TYR:CD1	1:D:210:VAL:HG22	2.44	0.53
1:A:374:TYR:O	1:A:378:VAL:HG23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:TYR:OH	2:C:401:E1U:O6P	2.21	0.52
1:A:139:SER:HA	1:A:169:TYR:OH	2.11	0.51
1:B:176:PRO:HG3	1:B:203:THR:HA	1.92	0.51
1:C:152:THR:HG22	1:C:173:VAL:HG13	1.93	0.50
1:A:283:VAL:C	1:A:284:LEU:HD12	2.32	0.49
1:D:135:ILE:HG13	1:D:163:VAL:HG12	1.94	0.49
1:D:270:ALA:HB1	1:D:311:ILE:HG23	1.95	0.49
1:A:284:LEU:HD13	1:A:313:TYR:HB2	1.94	0.49
1:C:139:SER:HA	1:C:169:TYR:OH	2.12	0.49
1:C:287:PRO:O	1:C:298:ARG:NH1	2.46	0.49
1:D:84:PHE:CD2	1:D:90:PHE:HA	2.49	0.48
1:B:306:ILE:HD11	1:B:313:TYR:CD1	2.49	0.48
1:B:324:MET:HB3	1:B:381:ALA:HB2	1.95	0.47
1:A:120:ASN:HB3	1:A:128:ASP:OD2	2.15	0.47
1:B:82:MET:HE1	3:B:402:MPD:H4	1.97	0.47
1:D:72:GLY:O	1:D:75:ASP:HB2	2.15	0.47
1:B:82:MET:HE3	3:B:402:MPD:H52	1.96	0.47
1:B:33:GLU:HA	1:B:66:HIS:O	2.15	0.46
1:C:36:CYS:HA	1:C:59:PRO:HB2	1.96	0.46
1:B:380:ASN:N	1:B:380:ASN:HD22	2.13	0.46
1:B:76:ILE:HG23	1:B:76:ILE:O	2.15	0.46
1:D:34:ALA:HB3	1:D:71:LEU:HD11	1.96	0.46
1:B:82:MET:CE	3:B:402:MPD:HM1	2.46	0.46
1:C:154:SER:HB3	1:C:182:PHE:CE1	2.51	0.46
1:B:34:ALA:HB3	1:B:71:LEU:HD11	1.97	0.46
1:A:257:ALA:HB3	1:A:365:VAL:HG22	1.97	0.45
1:C:299:VAL:HG21	1:C:315:LEU:HD21	1.99	0.45
1:B:81[B]:GLU:OE1	1:B:81[B]:GLU:HA	2.17	0.45
1:C:40:PRO:HG2	1:C:60:VAL:HG22	1.98	0.45
1:A:54:GLY:HA3	1:A:240:ARG:HD3	1.97	0.45
1:B:148:VAL:O	1:B:171:ALA:HA	2.17	0.45
1:A:222:VAL:O	1:A:223:ASN:C	2.55	0.44
1:C:119:TYR:CE2	3:C:402:MPD:H53	2.52	0.44
1:C:337:VAL:HG12	1:C:371:GLY:HA3	2.00	0.44
1:A:134:ILE:HG21	1:A:222:VAL:HB	2.00	0.43
1:C:189:LEU:HD11	1:D:328:VAL:HA	2.01	0.43
1:C:186:LEU:N	1:C:187:PRO:CD	2.81	0.43
1:C:152:THR:HG21	1:C:157:GLY:HA3	2.00	0.43
1:C:203:THR:OG1	1:C:225:TYR:OH	2.37	0.43
1:C:37:LEU:HD21	1:C:46:TYR:CD2	2.54	0.43
1:D:120:ASN:HB3	1:D:128:ASP:OD2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:LEU:HD23	1:D:331:ALA:HB1	1.99	0.43
1:D:129:ARG:N	1:D:130:PRO:HD2	2.34	0.43
1:A:133:GLU:CB	1:A:233:ALA:HB2	2.49	0.42
1:A:24:ASN:HB2	1:A:49:ALA:HB2	2.02	0.42
1:A:133:GLU:HB2	1:A:233:ALA:HB2	2.01	0.42
1:B:216:ASN:HD22	1:B:216:ASN:HA	1.72	0.42
1:C:152:THR:CG2	1:C:173:VAL:HG13	2.48	0.42
1:C:176:PRO:HG3	1:C:203:THR:HA	2.01	0.42
1:B:82:MET:CE	3:B:402:MPD:H31	2.49	0.42
1:B:95:LYS:HA	1:B:96:PRO:C	2.41	0.41
1:D:91:PHE:C	1:D:91:PHE:CD2	2.93	0.41
1:C:138:LEU:HD23	1:C:142:VAL:HB	2.02	0.41
1:B:152:THR:CG2	1:B:173:VAL:CG1	2.98	0.41
1:D:322:GLU:HB3	1:D:347:LYS:HG2	2.02	0.41
1:A:254:ARG:NE	1:A:254:ARG:HA	2.36	0.41
1:C:6:ILE:HD12	1:C:58:ILE:CG2	2.50	0.41
1:B:88:LEU:HA	1:B:88:LEU:HD13	1.93	0.41
1:A:197:ASP:HB3	1:A:206:LEU:HD21	2.03	0.41
1:A:324:MET:O	1:A:328:VAL:HG23	2.20	0.41
1:A:153:SER:O	1:A:154:SER:CB	2.68	0.41
1:B:13:LEU:HD23	1:B:137[B]:ARG:NE	2.35	0.41
1:A:248:ARG:CZ	1:C:2:ALA:HB2	2.51	0.41
1:D:110:VAL:HG13	1:D:363:TYR:CD1	2.56	0.41
1:D:340:PRO:HD2	1:D:374:TYR:CD2	2.56	0.41
1:A:110:VAL:HG13	1:A:363:TYR:CD1	2.56	0.40
1:B:147:LEU:HD23	1:B:219:PHE:HB3	2.03	0.40
1:D:288:ALA:CB	1:D:319:THR:HG22	2.51	0.40
1:B:152:THR:HG22	1:B:173:VAL:CG1	2.52	0.40
1:B:240[B]:ARG:HH11	1:B:240[B]:ARG:HD2	1.76	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	384/389 (99%)	365 (95%)	18 (5%)	1 (0%)	41	21
1	B	389/389 (100%)	372 (96%)	16 (4%)	1 (0%)	41	21
1	C	385/389 (99%)	370 (96%)	12 (3%)	3 (1%)	19	5
1	D	380/389 (98%)	366 (96%)	13 (3%)	1 (0%)	41	21
All	All	1538/1556 (99%)	1473 (96%)	59 (4%)	6 (0%)	34	15

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	154	SER
1	B	154	SER
1	D	154	SER
1	C	154	SER
1	C	39	ASN
1	C	38	GLY

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	312/312 (100%)	304 (97%)	8 (3%)	46	19
1	B	317/312 (102%)	303 (96%)	14 (4%)	28	7
1	C	313/312 (100%)	303 (97%)	10 (3%)	39	13
1	D	308/312 (99%)	295 (96%)	13 (4%)	30	7
All	All	1250/1248 (100%)	1205 (96%)	45 (4%)	36	10

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	CYS
1	A	88	LEU
1	A	101[A]	ARG
1	A	101[B]	ARG
1	A	104	LEU

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Mol	Chain	Res	Type
1	A	125	SER
1	A	203	THR
1	A	212	LYS
1	B	14	ASP
1	B	37	LEU
1	B	81[A]	GLU
1	B	81[B]	GLU
1	B	88	LEU
1	B	103	ARG
1	B	140	ARG
1	B	199	GLU
1	B	203	THR
1	B	212	LYS
1	B	216	ASN
1	B	236	ARG
1	B	284	LEU
1	B	309	LEU
1	C	20	SER
1	C	101	ARG
1	C	104	LEU
1	C	175	LEU
1	C	203	THR
1	C	226	TYR
1	C	248	ARG
1	C	314	THR
1	C	325	GLU
1	C	370	THR
1	D	13	LEU
1	D	35	ARG
1	D	75	ASP
1	D	103	ARG
1	D	104	LEU
1	D	138	LEU
1	D	203	THR
1	D	236	ARG
1	D	248	ARG
1	D	284	LEU
1	D	309	LEU
1	D	314	THR
1	D	383	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	380	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	MPD	B	402	-	7,7,7	0.38	0	9,10,10	0.80	0
2	E1U	B	401	-	23,26,26	1.98	8 (34%)	30,38,38	2.02	9 (30%)
2	E1U	D	401	-	23,26,26	2.02	7 (30%)	30,38,38	1.74	6 (20%)
2	E1U	A	401	-	23,26,26	2.09	8 (34%)	30,38,38	1.78	7 (23%)
2	E1U	C	401	-	23,26,26	1.96	7 (30%)	30,38,38	2.06	7 (23%)
3	MPD	C	402	-	7,7,7	0.32	0	9,10,10	0.67	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	MPD	B	402	-	-	0/5/5/5	-
2	E1U	B	401	-	-	5/17/21/21	0/1/1/1
2	E1U	D	401	-	-	5/17/21/21	0/1/1/1
2	E1U	A	401	-	-	5/17/21/21	0/1/1/1
2	E1U	C	401	-	-	5/17/21/21	0/1/1/1
3	MPD	C	402	-	-	1/5/5/5	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	E1U	C2A-C2	-5.55	1.41	1.50
2	D	401	E1U	C2A-C2	-5.14	1.41	1.50
2	B	401	E1U	C2A-C2	-5.04	1.41	1.50
2	C	401	E1U	C2A-C2	-4.96	1.42	1.50
2	C	401	E1U	C4A-N	3.94	1.34	1.27
2	D	401	E1U	C4-C4A	-3.77	1.39	1.46
2	A	401	E1U	P2-O6P	3.65	1.62	1.50
2	D	401	E1U	P2-O6P	3.47	1.61	1.50
2	A	401	E1U	C4A-N	3.28	1.33	1.27
2	B	401	E1U	P2-O6P	3.19	1.60	1.50
2	C	401	E1U	C5A-C5	-3.02	1.42	1.50
2	B	401	E1U	C4-C4A	-2.98	1.41	1.46
2	D	401	E1U	C4A-N	2.96	1.32	1.27
2	C	401	E1U	P2-O6P	2.95	1.60	1.50
2	B	401	E1U	C4A-N	2.93	1.32	1.27
2	A	401	E1U	C5A-C5	-2.87	1.43	1.50
2	B	401	E1U	C5A-C5	-2.84	1.43	1.50
2	D	401	E1U	C5A-C5	-2.81	1.43	1.50
2	C	401	E1U	C4-C4A	-2.79	1.41	1.46
2	A	401	E1U	C4-C4A	-2.66	1.41	1.46
2	B	401	E1U	C6-N1	2.59	1.39	1.34
2	C	401	E1U	C6-N1	2.49	1.39	1.34
2	A	401	E1U	C6-N1	2.47	1.39	1.34
2	B	401	E1U	P2-O5P	-2.46	1.45	1.54
2	D	401	E1U	P2-O5P	-2.44	1.45	1.54
2	D	401	E1U	C6-N1	2.44	1.39	1.34
2	C	401	E1U	P2-O5P	-2.36	1.45	1.54
2	A	401	E1U	P2-O5P	-2.19	1.46	1.54
2	A	401	E1U	P2-OG	2.18	1.67	1.60
2	B	401	E1U	P2-O7P	2.12	1.63	1.54

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	E1U	CA-N-C4A	6.70	126.49	117.40
2	B	401	E1U	CA-N-C4A	6.06	125.62	117.40
2	A	401	E1U	CA-N-C4A	5.03	124.22	117.40
2	D	401	E1U	CA-N-C4A	4.46	123.45	117.40
2	C	401	E1U	C4-C3-C2	4.37	122.89	120.19
2	A	401	E1U	C4-C3-C2	4.22	122.80	120.19
2	B	401	E1U	O4P-C5A-C5	4.22	117.39	109.35
2	D	401	E1U	O4P-C5A-C5	4.05	117.08	109.35
2	C	401	E1U	O4P-C5A-C5	3.81	116.60	109.35
2	C	401	E1U	O5P-P2-OG	3.50	116.04	106.73
2	B	401	E1U	C4-C3-C2	3.47	122.34	120.19
2	A	401	E1U	O4P-C5A-C5	3.36	115.75	109.35
2	B	401	E1U	C3-C2-N1	-3.02	116.86	120.77
2	D	401	E1U	CB-CA-N	-2.98	104.18	109.04
2	D	401	E1U	C4-C4A-N	-2.92	116.52	123.01
2	A	401	E1U	O5P-P2-OG	2.79	114.17	106.73
2	B	401	E1U	O5P-P2-OG	2.55	113.53	106.73
2	B	401	E1U	C2A-C2-C3	2.40	123.85	120.89
2	C	401	E1U	C3-C2-N1	-2.38	117.70	120.77
2	C	401	E1U	C2A-C2-C3	2.29	123.72	120.89
2	B	401	E1U	O7P-P2-O6P	-2.22	101.98	110.68
2	C	401	E1U	O7P-P2-O6P	-2.18	102.14	110.68
2	A	401	E1U	O7P-P2-O6P	-2.16	102.22	110.68
2	A	401	E1U	C3-C2-N1	-2.14	118.00	120.77
2	A	401	E1U	O2P-P-O3P	2.12	115.74	107.64
2	D	401	E1U	C3-C4-C5	2.10	119.87	118.26
2	B	401	E1U	C6-N1-C2	2.06	122.98	119.17
2	D	401	E1U	C3-C2-N1	-2.06	118.11	120.77
2	B	401	E1U	C4-C4A-N	-2.02	118.53	123.01

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	E1U	CB-OG-P2-O6P
2	B	401	E1U	C-CA-CB-OG
2	D	401	E1U	CB-OG-P2-O7P
2	D	401	E1U	C-CA-CB-OG
2	A	401	E1U	CB-OG-P2-O7P
2	A	401	E1U	CB-OG-P2-O5P
2	A	401	E1U	C-CA-CB-OG
2	C	401	E1U	C-CA-CB-OG
2	A	401	E1U	C-CA-N-C4A

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Mol	Chain	Res	Type	Atoms
2	A	401	E1U	CB-OG-P2-O6P
2	C	401	E1U	CB-OG-P2-O6P
2	C	401	E1U	CB-OG-P2-O5P
2	D	401	E1U	C3-C4-C4A-N
2	C	401	E1U	C3-C4-C4A-N
2	B	401	E1U	C5A-O4P-P-O1P
2	B	401	E1U	C-CA-N-C4A
3	C	402	MPD	O2-C2-C3-C4
2	D	401	E1U	CB-OG-P2-O5P
2	C	401	E1U	CB-OG-P2-O7P
2	B	401	E1U	C3-C4-C4A-N
2	D	401	E1U	C5-C4-C4A-N

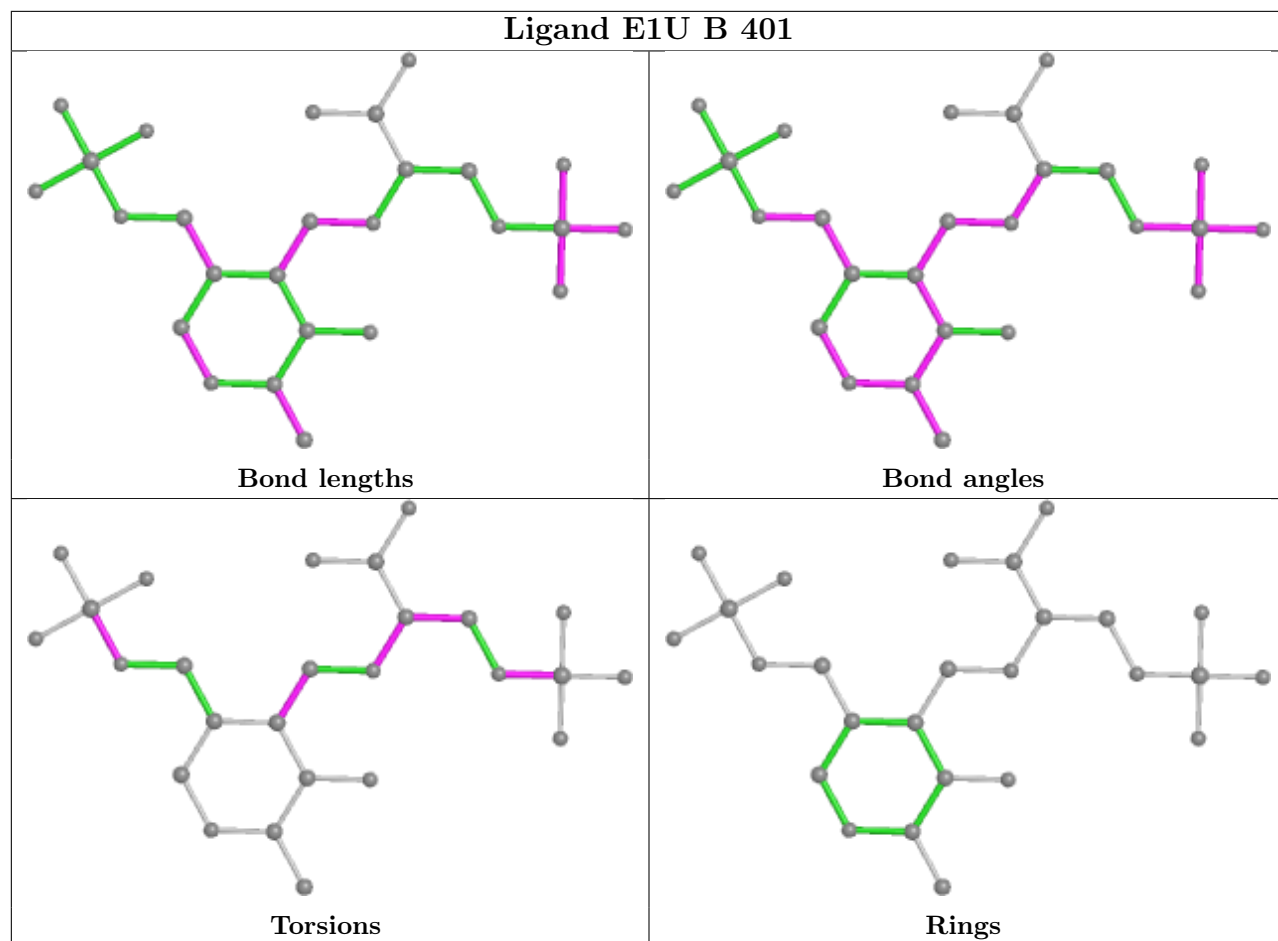
There are no ring outliers.

5 monomers are involved in 9 short contacts:

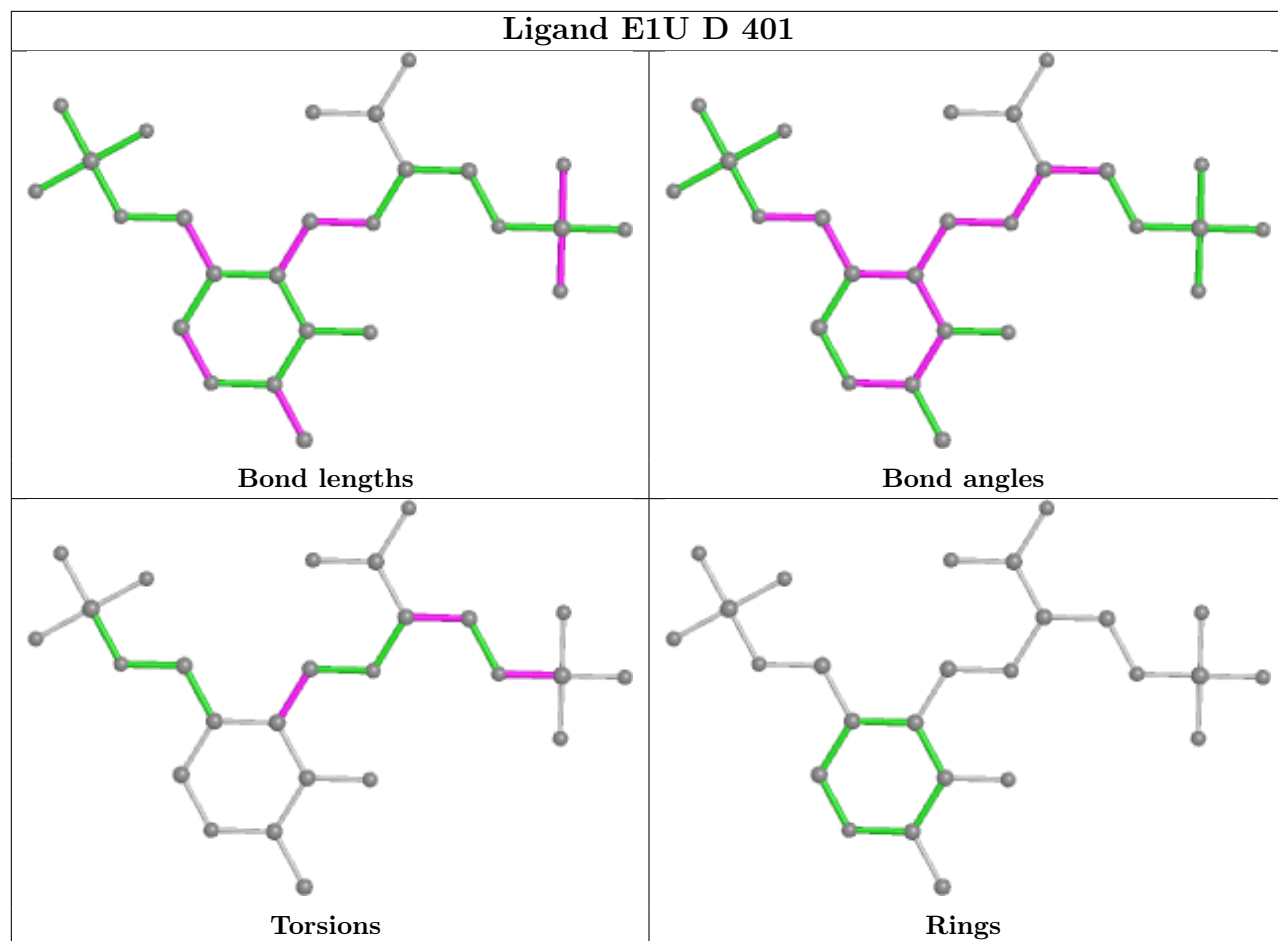
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	MPD	4	0
2	B	401	E1U	1	0
2	D	401	E1U	1	0
2	C	401	E1U	1	0
3	C	402	MPD	2	0

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.

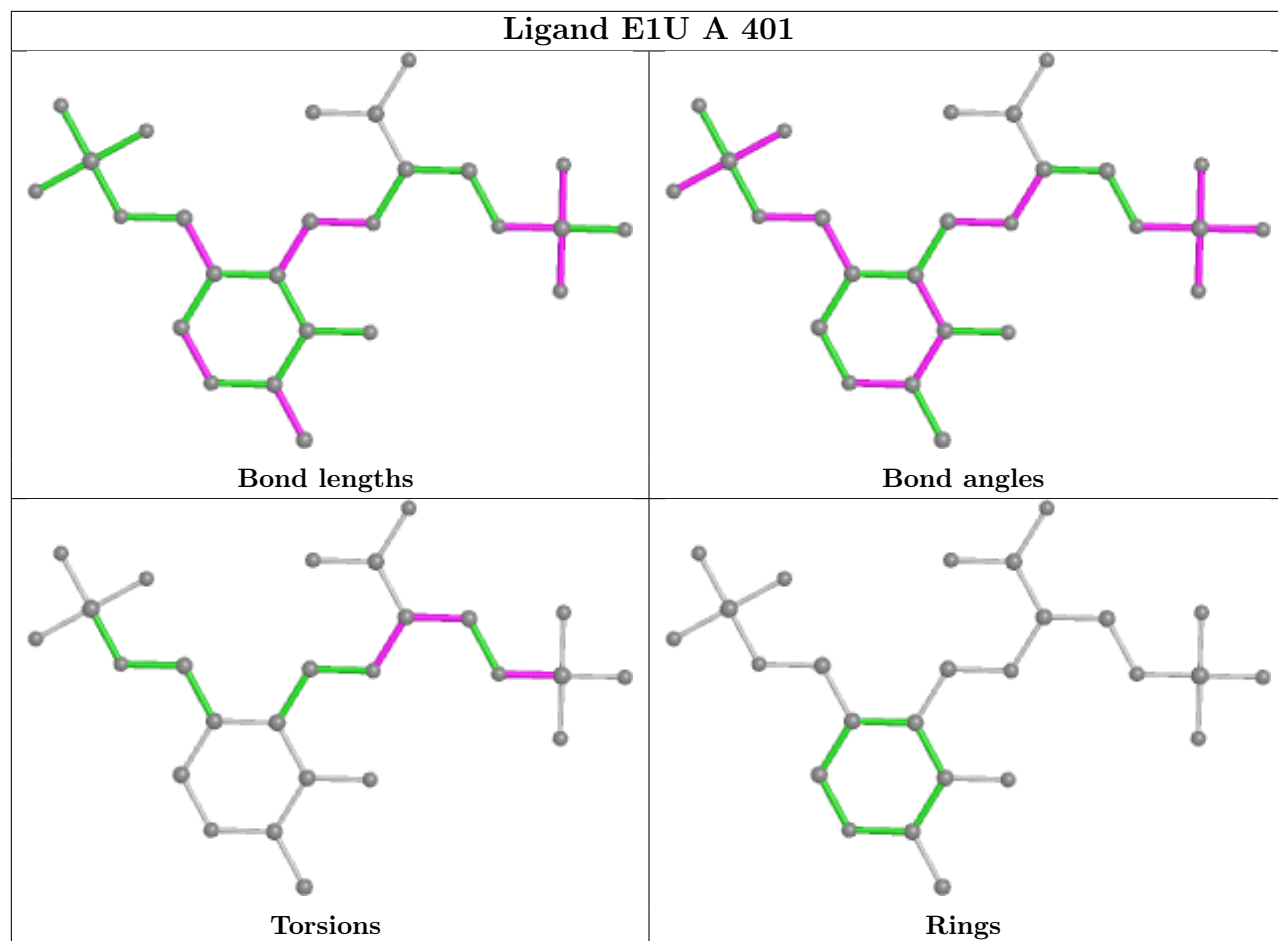
Ligand E1U B 401

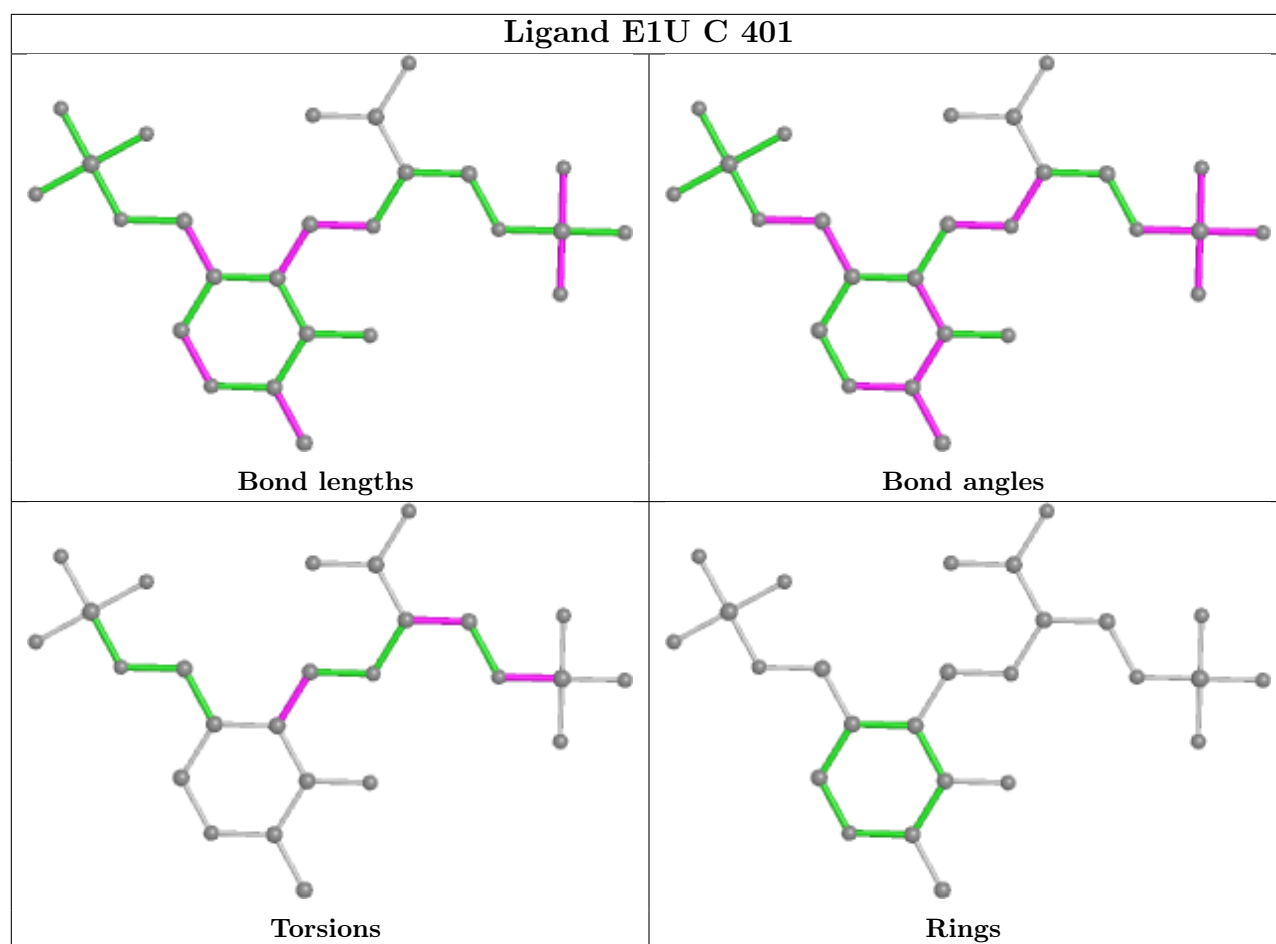


Ligand E1U D 401



Ligand E1U A 401





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	382/389 (98%)	0.09	15 (3%) 39 40	12, 24, 41, 63	0
1	B	382/389 (98%)	0.04	10 (2%) 56 57	11, 22, 39, 53	0
1	C	382/389 (98%)	0.08	15 (3%) 39 40	12, 23, 41, 65	0
1	D	382/389 (98%)	0.01	9 (2%) 59 60	10, 22, 39, 59	0
All	All	1528/1556 (98%)	0.05	49 (3%) 47 49	10, 23, 41, 65	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	37	LEU	11.8
1	C	62	ASP	10.3
1	B	37	LEU	9.7
1	A	2	ALA	6.7
1	C	37	LEU	5.7
1	A	63	GLY	5.0
1	A	62	ASP	4.9
1	A	14	ASP	4.5
1	C	38	GLY	4.4
1	A	38	GLY	4.2
1	C	383	GLU	4.1
1	D	35	ARG	3.7
1	D	38	GLY	3.7
1	B	64	CYS	3.7
1	A	39	ASN	3.7
1	C	63	GLY	3.6
1	A	383	GLU	3.6
1	A	37[A]	LEU	3.6
1	C	2	ALA	3.5
1	A	60	VAL	3.4
1	C	64	CYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	141	ARG	3.0
1	C	290	GLY	3.0
1	C	14	ASP	3.0
1	C	36	CYS	2.9
1	D	383	GLU	2.8
1	A	42	SER	2.7
1	B	290	GLY	2.7
1	D	205	HIS	2.6
1	B	36	CYS	2.6
1	C	205	HIS	2.6
1	D	140	ARG	2.5
1	B	199	GLU	2.5
1	A	40	PRO	2.4
1	B	60	VAL	2.4
1	D	302	GLY	2.3
1	A	140	ARG	2.3
1	B	39	ASN	2.3
1	B	61	GLY	2.2
1	C	289	GLN	2.2
1	A	64	CYS	2.2
1	B	76	ILE	2.2
1	A	302	GLY	2.1
1	B	141	ARG	2.1
1	A	306	ILE	2.1
1	C	199	GLU	2.1
1	D	64	CYS	2.1
1	C	60	VAL	2.0
1	D	307	ASN	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 monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

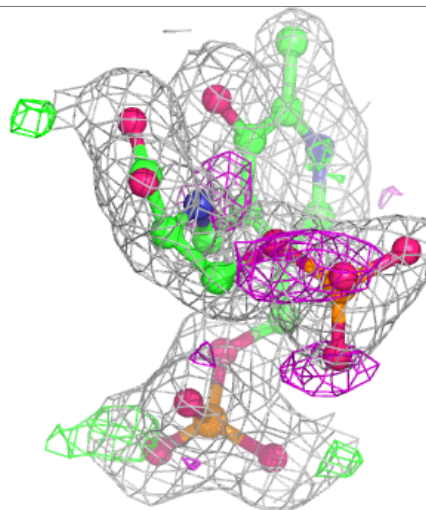
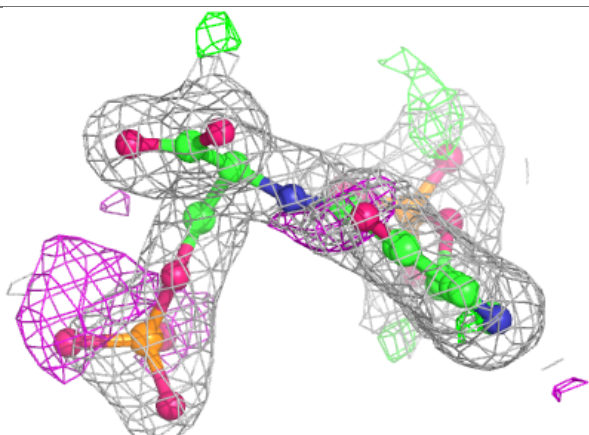
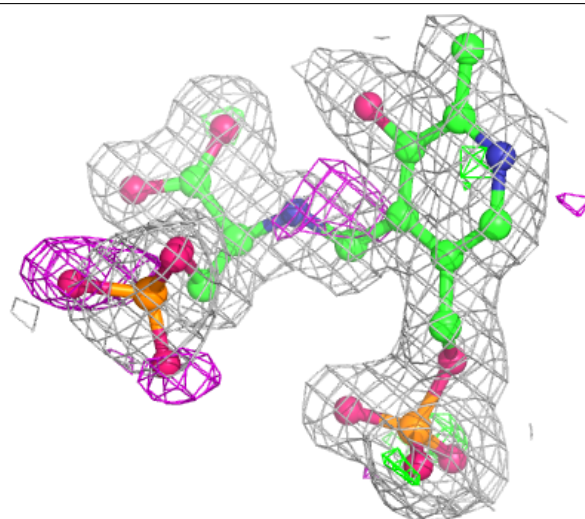
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MPD	C	402	8/8	0.77	0.16	41,42,43,43	0
3	MPD	B	402	8/8	0.82	0.12	39,39,40,41	0
2	E1U	A	401	26/26	0.93	0.12	22,25,36,39	0
2	E1U	C	401	26/26	0.93	0.10	22,24,38,40	0
2	E1U	B	401	26/26	0.93	0.10	20,21,37,42	0
2	E1U	D	401	26/26	0.94	0.11	20,23,40,43	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.

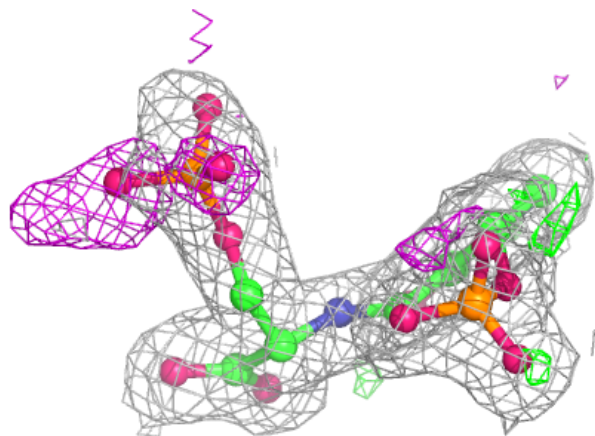
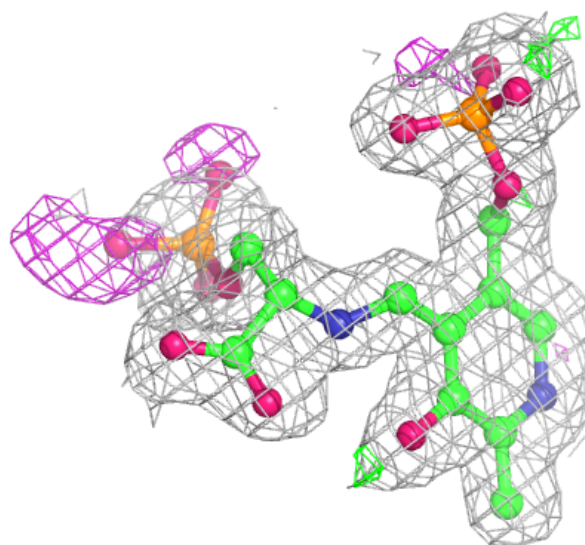
Electron density around E1U A 401:

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



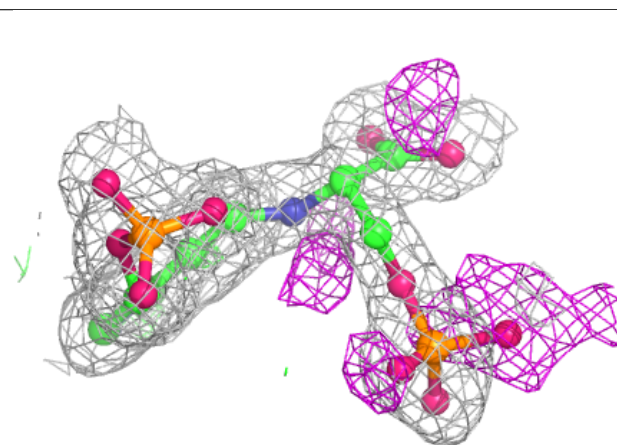
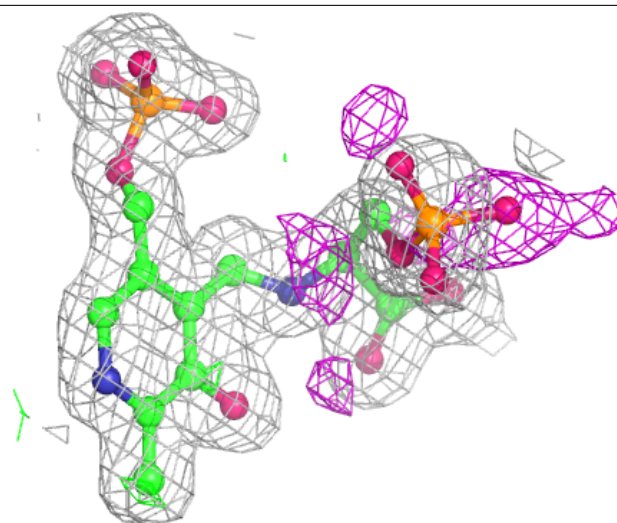
Electron density around E1U C 401:

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



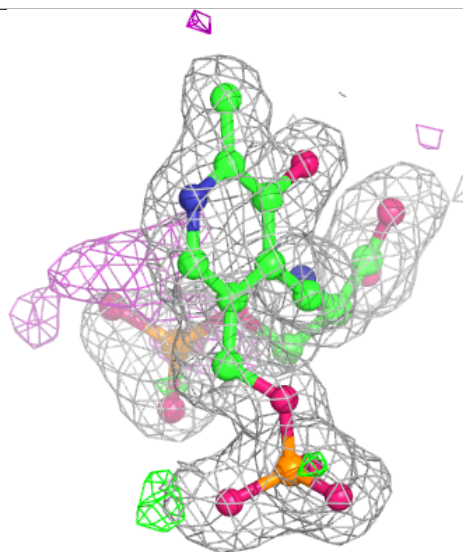
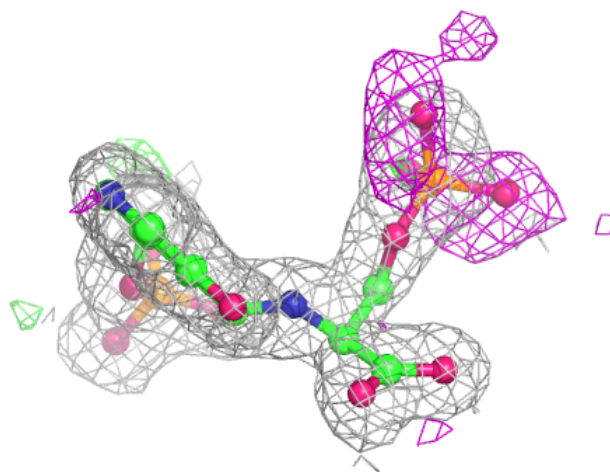
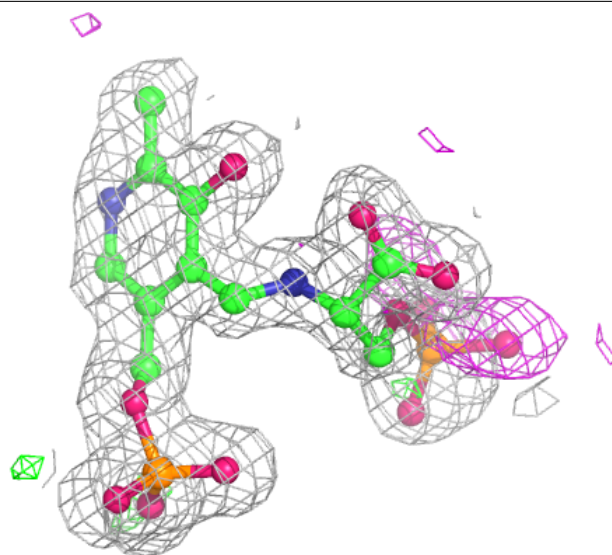
Electron density around E1U B 401:

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



Electron density around E1U D 401:

$2mF_o-DF_c$ (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.