



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:46 PM UTC

PDB ID : 5EUE / pdb_00005eue
Title : S1P Lyase Bacterial Surrogate bound to N-(2-((4-methoxy-2,5-dimethylbenzyl)amino)-1-phenylethyl)-5-methylisoxazole-3-carboxamide
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Deposited on : 2015-11-18
Resolution : 2.83 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](#) ①) 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

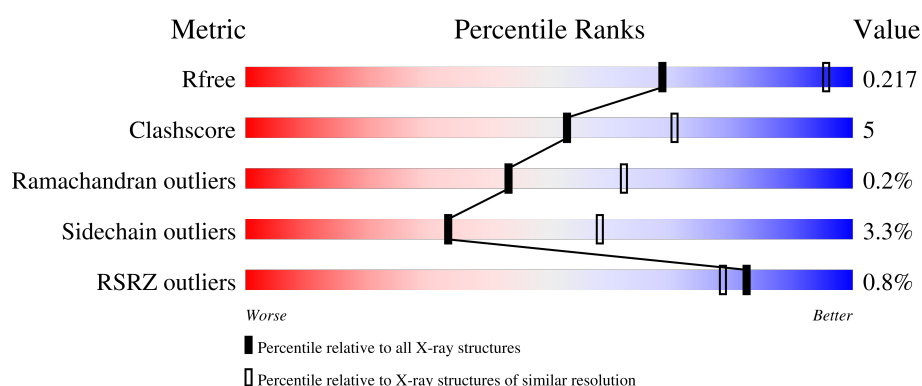
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.83 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	514	74% 13% 12%
1	B	514	74% 13% 13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7026 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.

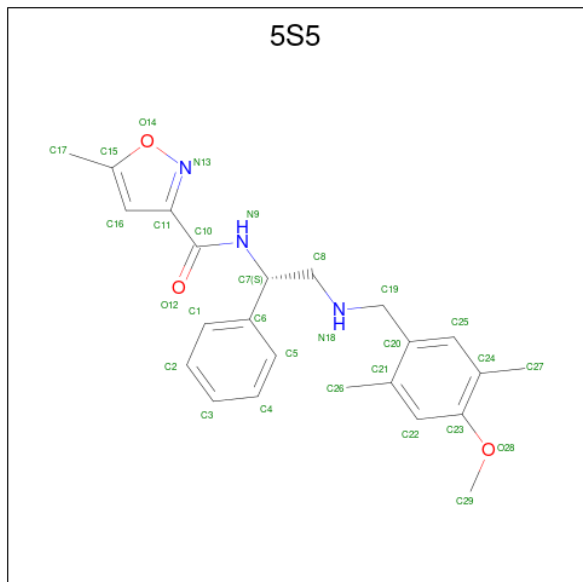
- Molecule 1 is a protein called Putative sphingosine-1-phosphate lyase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	P	S	0	0	0
			3423	2187	598	624	1	13			
1	B	449	Total	C	N	O	P	S	0	0	0
			3409	2176	597	622	1	13			

There are 24 discrepancies between the modelled and reference sequences:

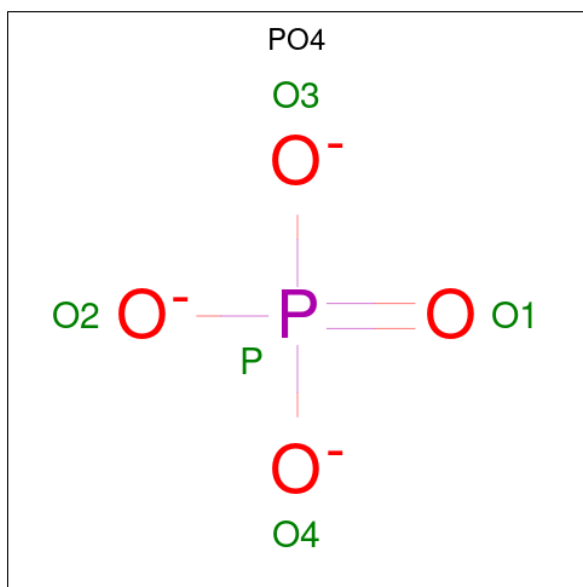
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q67PY4
A	1	PRO	-	expression tag	UNP Q67PY4
A	249	PHE	TYR	engineered mutation	UNP Q67PY4
A	344	ILE	LEU	engineered mutation	UNP Q67PY4
A	346	ALA	PHE	engineered mutation	UNP Q67PY4
A	497	SER	LEU	engineered mutation	UNP Q67PY4
A	508	HIS	-	expression tag	UNP Q67PY4
A	509	HIS	-	expression tag	UNP Q67PY4
A	510	HIS	-	expression tag	UNP Q67PY4
A	511	HIS	-	expression tag	UNP Q67PY4
A	512	HIS	-	expression tag	UNP Q67PY4
A	513	HIS	-	expression tag	UNP Q67PY4
B	0	MET	-	initiating methionine	UNP Q67PY4
B	1	PRO	-	expression tag	UNP Q67PY4
B	249	PHE	TYR	engineered mutation	UNP Q67PY4
B	344	ILE	LEU	engineered mutation	UNP Q67PY4
B	346	ALA	PHE	engineered mutation	UNP Q67PY4
B	497	SER	LEU	engineered mutation	UNP Q67PY4
B	508	HIS	-	expression tag	UNP Q67PY4
B	509	HIS	-	expression tag	UNP Q67PY4
B	510	HIS	-	expression tag	UNP Q67PY4
B	511	HIS	-	expression tag	UNP Q67PY4
B	512	HIS	-	expression tag	UNP Q67PY4
B	513	HIS	-	expression tag	UNP Q67PY4

- Molecule 2 is {N}-[(1 {S})-2-[(4-methoxy-2,5-dimethyl-phenyl)methylamino]-1-phenyl-ethyl]-5-methyl-1,2-oxazole-3-carboxamide (CCD ID: 5S5) (formula: $C_{23}H_{27}N_3O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			29	23	3	3		
2	B	1	Total	C	N	O	0	0
			29	23	3	3		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

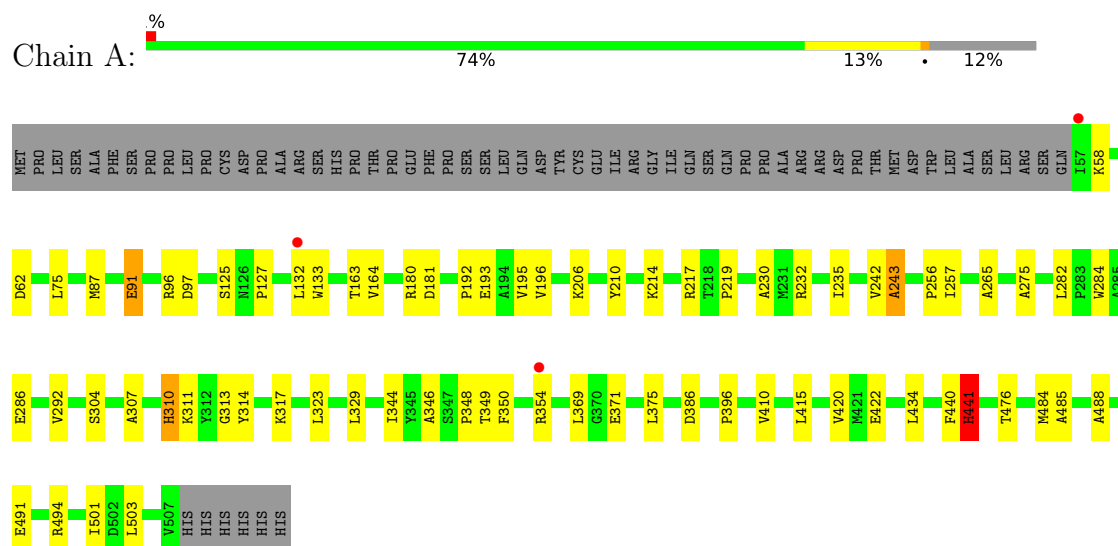
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total	O	0	0
			58	58		
4	B	68	Total	O	0	0
			68	68		

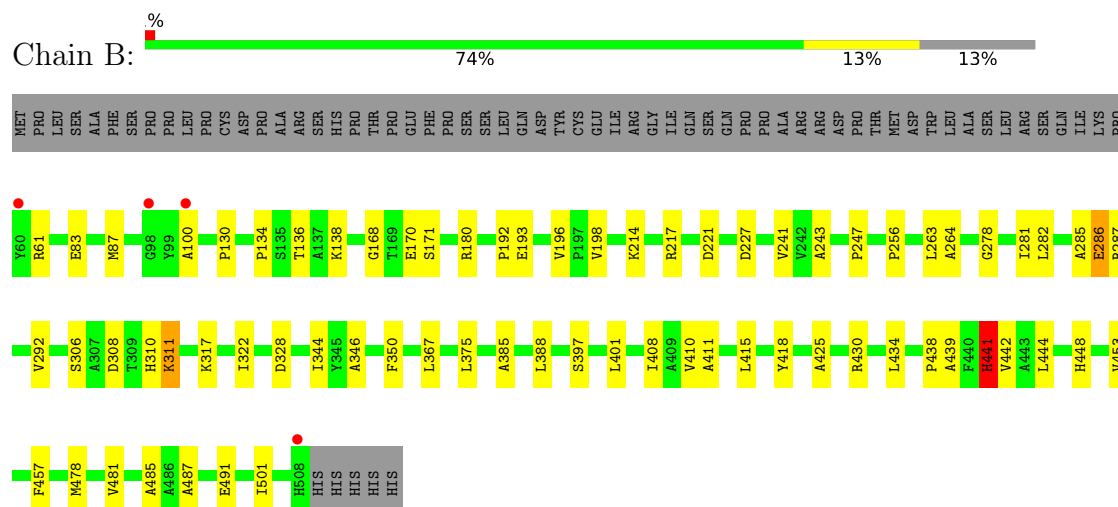
3 Residue-property plots

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: Putative sphingosine-1-phosphate lyase



- Molecule 1: Putative sphingosine-1-phosphate lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.01Å 129.61Å 84.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.09 – 2.83 71.09 – 2.83	Depositor EDS
% Data completeness (in resolution range)	99.6 (71.09-2.83) 99.8 (71.09-2.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.81Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.173 , 0.249 (Not available) , 0.217	Depositor DCC
R_{free} test set	1168 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	48.7	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7026	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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: PO4, LLP, 5S5

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	0.90	1/3494 (0.0%)	1.40	17/4773 (0.4%)
1	B	0.89	0/3480	1.39	16/4754 (0.3%)
All	All	0.90	1/6974 (0.0%)	1.40	33/9527 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	257	ILE	CA-CB	7.86	1.58	1.54

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	441	HIS	CA-CB-CG	6.43	120.23	113.80
1	A	441	HIS	CA-CB-CG	6.37	120.17	113.80
1	A	256	PRO	CA-C-N	6.37	126.54	120.43
1	A	256	PRO	C-N-CA	6.37	126.54	120.43
1	A	243	ALA	CA-C-N	6.20	127.35	121.46
1	A	243	ALA	C-N-CA	6.20	127.35	121.46
1	A	91	GLU	N-CA-C	5.91	118.95	111.69
1	A	386	ASP	CA-C-N	5.83	128.56	120.29
1	A	386	ASP	C-N-CA	5.83	128.56	120.29
1	A	313	GLY	N-CA-C	-5.80	107.46	114.66
1	B	256	PRO	CA-C-N	5.76	123.87	120.24
1	B	256	PRO	C-N-CA	5.76	123.87	120.24
1	B	285	ALA	CA-C-N	5.71	128.19	120.65
1	B	285	ALA	C-N-CA	5.71	128.19	120.65
1	A	265	ALA	CA-C-N	5.68	127.89	120.28
1	A	265	ALA	C-N-CA	5.68	127.89	120.28
1	B	487	ALA	CA-C-N	5.63	129.22	120.60
1	B	487	ALA	C-N-CA	5.63	129.22	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	GLY	CA-C-N	5.61	132.06	123.47
1	A	313	GLY	C-N-CA	5.61	132.06	123.47
1	B	227	ASP	CA-CB-CG	5.54	118.14	112.60
1	A	181	ASP	CA-C-N	5.46	127.54	120.44
1	A	181	ASP	C-N-CA	5.46	127.54	120.44
1	B	241	VAL	N-CA-C	5.44	115.39	107.88
1	B	221	ASP	CA-CB-CG	5.26	117.86	112.60
1	A	386	ASP	CA-CB-CG	5.25	117.85	112.60
1	B	130	PRO	CA-C-N	5.19	127.75	120.28
1	B	130	PRO	C-N-CA	5.19	127.75	120.28
1	A	62	ASP	CA-CB-CG	5.12	117.72	112.60
1	B	264	ALA	CA-C-N	5.01	127.00	120.28
1	B	264	ALA	C-N-CA	5.01	127.00	120.28
1	B	425	ALA	CA-C-N	5.01	125.54	119.98
1	B	425	ALA	C-N-CA	5.01	125.54	119.98

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	3423	0	3357	38	0
1	B	3409	0	3333	38	0
2	A	29	0	0	0	0
2	B	29	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	58	0	0	0	0
4	B	68	0	0	0	0
All	All	7026	0	6690	68	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 (68) 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:286:GLU:HG3	1:A:292:VAL:HG13	1.58	0.85
1:B:196:VAL:HG12	1:B:243:ALA:HB3	1.60	0.83
1:B:385:ALA:HB2	1:B:444:LEU:HD12	1.68	0.74
1:A:284:TRP:HB3	1:A:375:LEU:HD23	1.69	0.73
1:B:286:GLU:HG3	1:B:292:VAL:CG1	2.17	0.73
1:A:196:VAL:HG12	1:A:243:ALA:HB3	1.70	0.72
1:B:434:LEU:HD11	1:B:441:HIS:HB3	1.70	0.71
1:A:415:LEU:HD11	1:A:420:VAL:HG21	1.73	0.70
1:B:282:LEU:HD22	1:B:292:VAL:HG21	1.74	0.69
1:B:286:GLU:HG3	1:B:292:VAL:HG12	1.75	0.67
1:B:388:LEU:HD13	1:B:444:LEU:HD21	1.79	0.65
1:B:310:HIS:HB2	1:B:317:LYS:HA	1.78	0.64
1:A:344:ILE:HD11	1:B:485:ALA:HB1	1.83	0.60
1:A:282:LEU:HD22	1:A:292:VAL:HG21	1.83	0.58
1:A:501:ILE:HD11	1:B:346:ALA:HB1	1.86	0.58
1:A:346:ALA:HB1	1:B:501:ILE:HD11	1.86	0.57
1:A:127:PRO:HG2	1:A:354:ARG:HG2	1.87	0.57
1:A:485:ALA:HB1	1:B:344:ILE:HD11	1.85	0.57
1:A:193:GLU:HG2	1:A:214:LYS:HB3	1.88	0.56
1:A:415:LEU:CD1	1:A:420:VAL:CG2	2.84	0.56
1:A:195:VAL:HB	1:A:242:VAL:HG12	1.88	0.55
1:A:434:LEU:HD11	1:A:441:HIS:HB3	1.87	0.55
1:A:91:GLU:OE2	1:B:134:PRO:HD2	2.06	0.55
1:A:219:PRO:HD2	1:A:230:ALA:HB1	1.89	0.54
1:B:193:GLU:HG2	1:B:214:LYS:HB3	1.90	0.54
1:B:478:MET:HE3	1:B:481:VAL:HG11	1.89	0.54
1:A:415:LEU:HD11	1:A:420:VAL:CG2	2.36	0.54
1:B:83:GLU:HG2	1:B:87:MET:HE3	1.91	0.53
1:A:410:VAL:CG1	1:A:440:PHE:CZ	2.92	0.52
1:B:170:GLU:HG3	1:B:350:PHE:CD2	2.46	0.50
1:A:371:GLU:O	1:A:375:LEU:HG	2.11	0.50
1:B:180:ARG:HG3	1:B:192:PRO:HG3	1.93	0.50
1:A:180:ARG:HG3	1:A:192:PRO:HG3	1.94	0.48
1:A:310:HIS:HB2	1:A:317:LYS:HA	1.95	0.48
1:B:308:ASP:HB3	1:B:310:HIS:HD2	1.79	0.48
1:A:96:ARG:HD2	1:A:97:ASP:OD2	2.13	0.48
1:B:308:ASP:HB3	1:B:310:HIS:CD2	2.48	0.48
1:A:164:VAL:HG12	1:A:354:ARG:HH21	1.78	0.47
1:A:232:ARG:HA	1:A:235:ILE:HD12	1.96	0.47
1:B:247:PRO:HD3	1:B:278:GLY:HA3	1.97	0.47
1:A:163:THR:HG21	1:A:348:PRO:HB2	1.97	0.46
1:A:314:TYR:HD2	1:A:369:LEU:CD1	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:418:TYR:CE2	1:B:438:PRO:HB3	2.50	0.46
1:B:411:ALA:HB2	1:B:439:ALA:HB2	1.98	0.46
1:A:206:LYS:HG2	1:A:210:TYR:CZ	2.51	0.46
1:B:441:HIS:CD2	1:B:441:HIS:C	2.95	0.45
1:A:422:GLU:HG2	1:A:476:THR:O	2.17	0.44
1:A:125:SER:HA	1:A:133:TRP:CE2	2.52	0.44
1:B:388:LEU:HD22	1:B:408:ILE:HG13	1.99	0.44
1:B:448:HIS:HA	1:B:453:VAL:HG11	2.00	0.44
1:B:198:VAL:HG12	1:B:217:ARG:HB3	1.98	0.44
1:B:385:ALA:CB	1:B:444:LEU:HD12	2.41	0.44
1:A:132:LEU:HD21	1:B:430:ARG:HG3	2.00	0.43
1:B:442:VAL:HG21	1:B:457:PHE:CZ	2.54	0.42
1:A:196:VAL:O	1:A:217:ARG:HA	2.19	0.42
1:A:304:SER:HA	1:A:323:LEU:O	2.19	0.42
1:A:484:MET:O	1:A:488:ALA:HB2	2.19	0.42
1:B:401:LEU:HD21	1:B:411:ALA:HB3	2.02	0.41
1:A:275:ALA:O	1:A:307:ALA:HA	2.20	0.41
1:B:168:GLY:HA3	1:B:311:LLP:C5'	2.50	0.41
1:A:75:LEU:HD12	1:B:367:LEU:HB3	2.02	0.41
1:A:282:LEU:HD22	1:A:292:VAL:CG2	2.49	0.41
1:B:100:ALA:HA	1:B:430:ARG:HB2	2.03	0.41
1:A:441:HIS:C	1:A:441:HIS:CD2	2.99	0.40
1:B:171:SER:HB3	1:B:306:SER:HB2	2.03	0.40
1:B:317:LYS:HA	1:B:317:LYS:HD3	1.83	0.40
1:B:385:ALA:HB2	1:B:444:LEU:CD1	2.44	0.40
1:A:87:MET:HE1	1:B:138:LYS:HE3	2.03	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	448/514 (87%)	427 (95%)	20 (4%)	1 (0%)	43	62
1	B	446/514 (87%)	431 (97%)	14 (3%)	1 (0%)	43	62
All	All	894/1028 (87%)	858 (96%)	34 (4%)	2 (0%)	43	62

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	61	ARG
1	A	396	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	336/394 (85%)	327 (97%)	9 (3%)	39	63
1	B	334/394 (85%)	321 (96%)	13 (4%)	28	53
All	All	670/788 (85%)	648 (97%)	22 (3%)	33	58

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	LYS
1	A	310	HIS
1	A	329	LEU
1	A	349	THR
1	A	350	PHE
1	A	441	HIS
1	A	491	GLU
1	A	494	ARG
1	A	503	LEU
1	B	136	THR
1	B	263	LEU
1	B	281	ILE
1	B	286	GLU
1	B	287	ARG

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Mol	Chain	Res	Type
1	B	322	ILE
1	B	328	ASP
1	B	375	LEU
1	B	397	SER
1	B	410	VAL
1	B	415	LEU
1	B	441	HIS
1	B	491	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	116	ASN
1	A	122	GLN
1	A	383	GLN
1	A	441	HIS
1	A	467	GLN
1	B	106	HIS
1	B	122	GLN
1	B	333	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	LLP	B	311	1	23,24,25	2.58	7 (30%)	25,32,34	2.34	8 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	LLP	A	311	1	23,24,25	2.73	8 (34%)	25,32,34	1.81	9 (36%)

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
1	LLP	B	311	1	-	8/16/17/19	0/1/1/1
1	LLP	A	311	1	-	7/16/17/19	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	311	LLP	C4'-NZ	8.03	1.54	1.27
1	B	311	LLP	C4'-NZ	6.07	1.47	1.27
1	B	311	LLP	C3-C2	6.04	1.47	1.41
1	B	311	LLP	C4-C4'	5.39	1.58	1.46
1	A	311	LLP	C4-C4'	4.72	1.56	1.46
1	A	311	LLP	C4-C5	4.50	1.48	1.42
1	A	311	LLP	C3-C2	4.24	1.45	1.41
1	A	311	LLP	P-OP4	-4.14	1.47	1.60
1	B	311	LLP	C4-C3	3.52	1.46	1.41
1	B	311	LLP	C4-C5	3.49	1.46	1.42
1	B	311	LLP	CB-CA	3.05	1.58	1.53
1	A	311	LLP	C6-C5	2.81	1.43	1.37
1	B	311	LLP	P-OP4	-2.52	1.52	1.60
1	A	311	LLP	C2-N1	2.11	1.37	1.33
1	A	311	LLP	C2'-C2	2.08	1.53	1.50

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	311	LLP	OP3-P-OP4	5.26	120.38	106.67
1	B	311	LLP	C5'-C5-C6	-4.83	111.49	119.36
1	A	311	LLP	C4-C4'-NZ	-4.19	104.69	124.04
1	B	311	LLP	C3-C4-C5	-4.13	114.96	118.28
1	B	311	LLP	C2'-C2-C3	3.97	125.45	120.80
1	B	311	LLP	OP4-C5'-C5	3.34	115.62	109.36
1	A	311	LLP	C4-C3-C2	-2.99	118.46	120.14
1	B	311	LLP	OP4-P-OP1	-2.90	98.60	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	311	LLP	CD-CE-NZ	-2.84	103.31	110.83
1	A	311	LLP	O3-C3-C2	2.79	123.37	117.58
1	A	311	LLP	C5-C4-C4'	2.77	125.75	121.47
1	A	311	LLP	C5-C6-N1	-2.57	119.65	123.83
1	B	311	LLP	CE-NZ-C4'	-2.50	110.71	118.72
1	A	311	LLP	C5'-C5-C6	-2.40	115.45	119.36
1	A	311	LLP	C3-C4-C4'	-2.33	116.20	120.40
1	B	311	LLP	C6-N1-C2	2.26	123.30	119.20
1	A	311	LLP	C6-N1-C2	2.24	123.26	119.20

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	311	LLP	C4-C4'-NZ-CE
1	A	311	LLP	C4-C5-C5'-OP4
1	A	311	LLP	C6-C5-C5'-OP4
1	A	311	LLP	O-C-CA-CB
1	B	311	LLP	C4-C4'-NZ-CE
1	B	311	LLP	C5'-OP4-P-OP2
1	B	311	LLP	O-C-CA-CB
1	A	311	LLP	C3-C4-C4'-NZ
1	A	311	LLP	C5-C4-C4'-NZ
1	B	311	LLP	C3-C4-C4'-NZ
1	B	311	LLP	C5'-OP4-P-OP3
1	B	311	LLP	C5'-OP4-P-OP1
1	B	311	LLP	CG-CD-CE-NZ
1	A	311	LLP	CD-CE-NZ-C4'
1	B	311	LLP	C5-C4-C4'-NZ

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	311	LLP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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	PO4	A	602	-	4,4,4	0.96	0	6,6,6	0.91	0
3	PO4	B	601	-	4,4,4	0.95	0	6,6,6	0.63	0
2	5S5	B	602	-	31,31,31	1.02	1 (3%)	39,42,42	1.96	5 (12%)
2	5S5	A	601	-	31,31,31	1.03	3 (9%)	39,42,42	1.42	3 (7%)

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
2	5S5	B	602	-	-	8/20/20/20	0/3/3/3
2	5S5	A	601	-	-	9/20/20/20	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	602	5S5	C11-C10	-2.57	1.43	1.48
2	A	601	5S5	C11-C10	-2.45	1.44	1.48
2	A	601	5S5	C10-N9	2.16	1.38	1.34
2	A	601	5S5	C16-C15	2.02	1.38	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	5S5	C29-O28-C23	-8.70	104.76	117.51
2	A	601	5S5	C29-O28-C23	-5.16	109.95	117.51
2	B	602	5S5	O28-C23-C24	3.66	121.43	115.09
2	B	602	5S5	O28-C23-C22	-3.66	117.77	124.08
2	A	601	5S5	O28-C23-C24	3.24	120.69	115.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	5S5	O28-C23-C22	-2.84	119.19	124.08
2	B	602	5S5	O14-C15-C17	2.12	118.80	115.89
2	B	602	5S5	C19-C20-C25	-2.03	116.43	120.34

There are no chirality outliers.

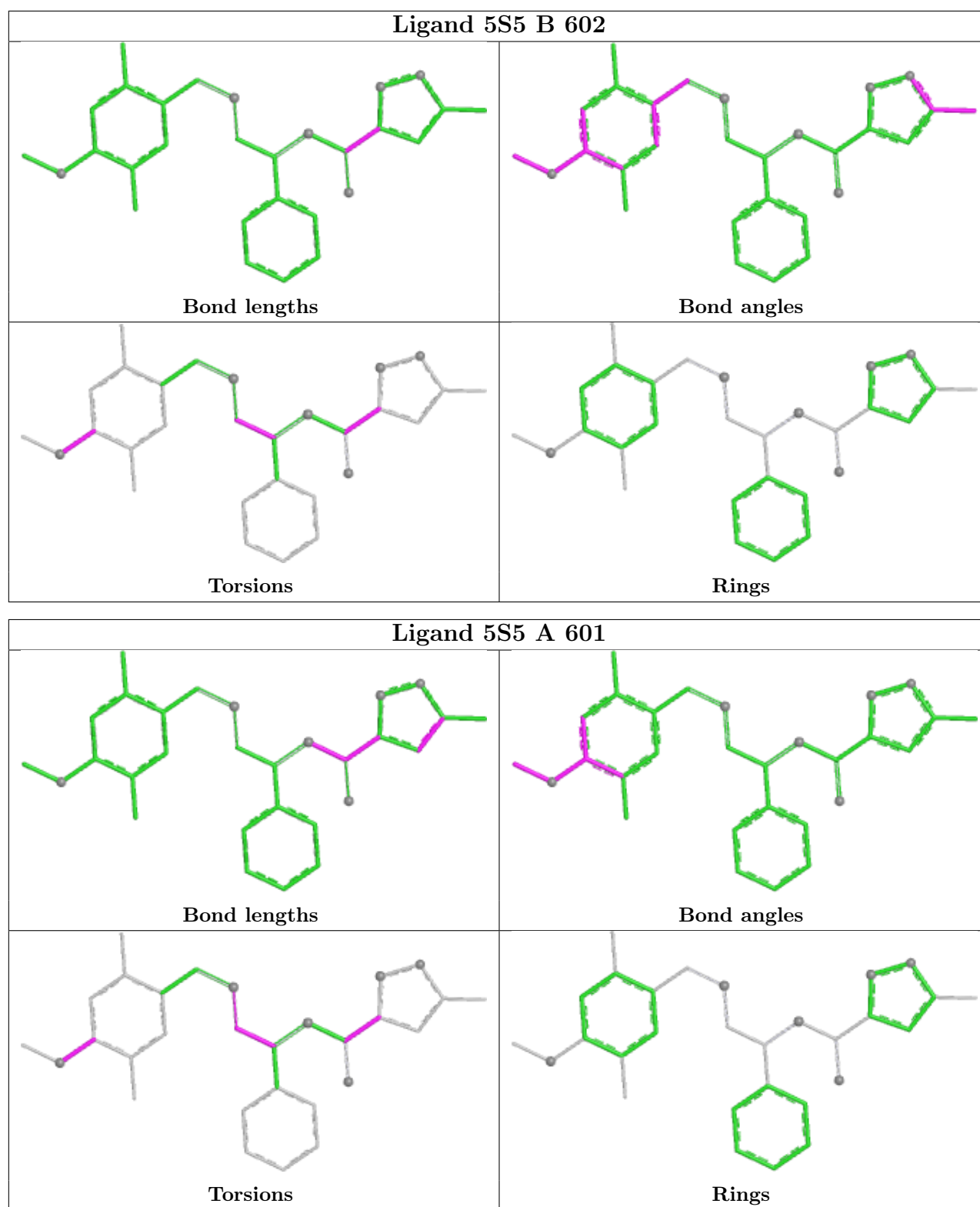
All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	5S5	C6-C7-C8-N18
2	A	601	5S5	N9-C7-C8-N18
2	A	601	5S5	N9-C10-C11-C16
2	A	601	5S5	N9-C10-C11-N13
2	A	601	5S5	O12-C10-C11-C16
2	A	601	5S5	O12-C10-C11-N13
2	B	602	5S5	C6-C7-C8-N18
2	B	602	5S5	N9-C10-C11-C16
2	B	602	5S5	O12-C10-C11-C16
2	A	601	5S5	C22-C23-O28-C29
2	B	602	5S5	C22-C23-O28-C29
2	B	602	5S5	C24-C23-O28-C29
2	A	601	5S5	C24-C23-O28-C29
2	B	602	5S5	N9-C7-C8-N18
2	B	602	5S5	N9-C10-C11-N13
2	B	602	5S5	O12-C10-C11-N13
2	A	601	5S5	C7-C8-N18-C19

There are no ring outliers.

No monomer is involved in short contacts.

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	450/514 (87%)	-0.24	3 (0%) 84 80	25, 41, 72, 100	0
1	B	448/514 (87%)	-0.23	4 (0%) 81 76	25, 42, 74, 120	0
All	All	898/1028 (87%)	-0.24	7 (0%) 82 78	25, 42, 73, 120	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	508	HIS	3.7
1	B	98	GLY	3.0
1	A	132	LEU	2.4
1	B	60	TYR	2.3
1	B	100	ALA	2.1
1	A	354	ARG	2.1
1	A	57	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	LLP	B	311	24/25	0.85	0.13	61,79,86,87	0
1	LLP	A	311	24/25	0.96	0.09	32,46,62,62	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

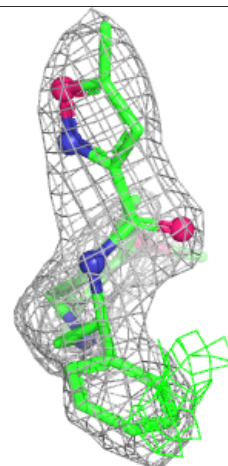
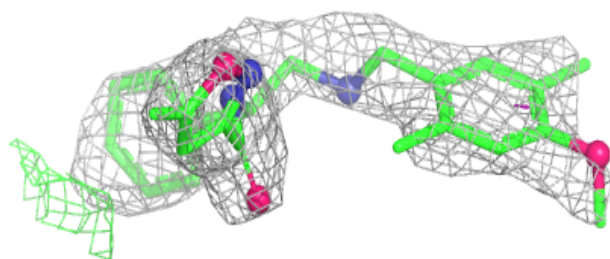
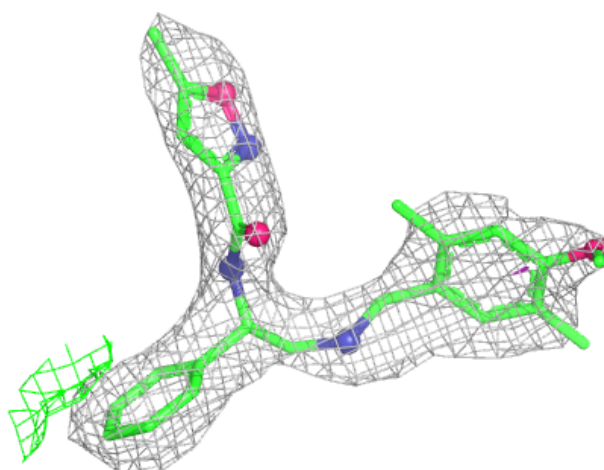
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	5S5	A	601	29/29	0.83	0.17	70,75,79,81	0
3	PO4	B	601	5/5	0.86	0.12	61,62,63,63	5
2	5S5	B	602	29/29	0.87	0.14	59,65,79,81	0
3	PO4	A	602	5/5	0.95	0.11	25,30,32,35	5

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.

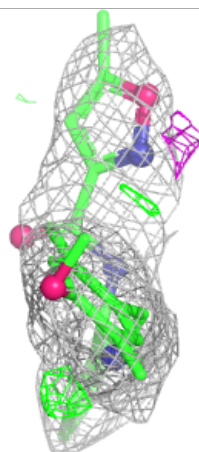
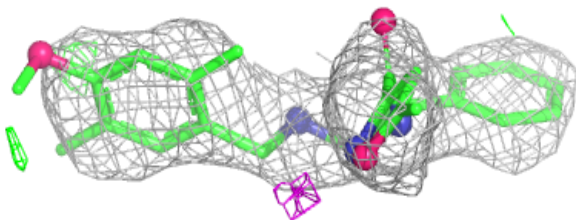
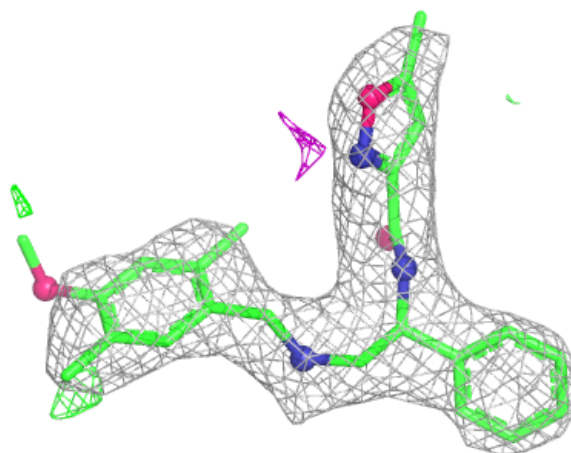
Electron density around 5S5 A 601:

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 5S5 B 602:

$2mF_o-DF_c$ (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.