



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 06:07 am BST

PDB ID : 4QDE
Title : Dcps in complex with covalent inhibitor
Authors : Liu, S.
Deposited on : 2014-05-13
Resolution : 2.90 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 : 2.11
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) : 2.11

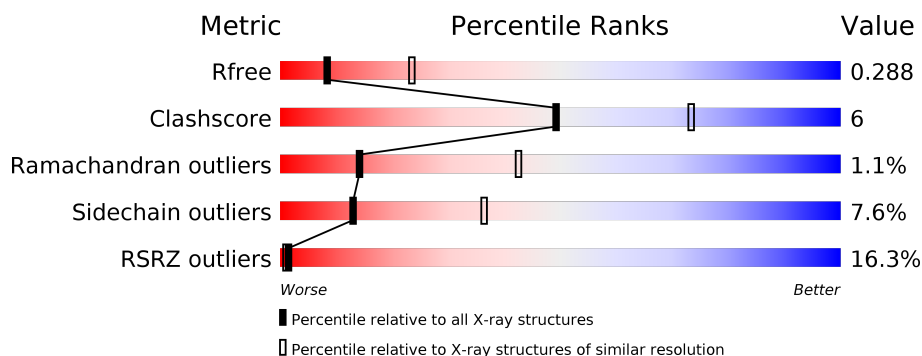
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 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	337	<div> <div>11%</div> <div> <div></div> <div>66%</div> <div>18%</div> <div>•</div> <div>15%</div> </div> </div>
1	B	337	<div> <div>14%</div> <div> <div></div> <div>68%</div> <div>16%</div> <div>•</div> <div>15%</div> </div> </div>
1	C	337	<div> <div>17%</div> <div> <div></div> <div>69%</div> <div>15%</div> <div>•</div> <div>14%</div> </div> </div>
1	D	337	<div> <div>14%</div> <div> <div></div> <div>69%</div> <div>16%</div> <div>•</div> <div>13%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	B	402	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9655 atoms, of which 13 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 m7GpppX diphosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2377	1523	425	426	3			
1	B	288	Total	C	N	O	S	0	0	0
			2384	1526	427	428	3			
1	C	289	Total	C	N	O	S	0	0	0
			2389	1529	427	430	3			
1	D	292	Total	C	N	O	S	0	0	0
			2411	1542	430	436	3			

There are 4 discrepancies between the modelled and reference sequences:

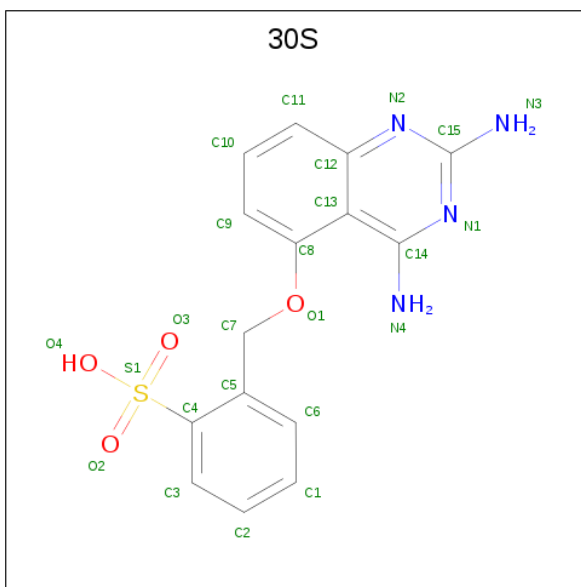
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	EXPRESSION TAG	UNP Q96C86
B	1	GLY	-	EXPRESSION TAG	UNP Q96C86
C	1	GLY	-	EXPRESSION TAG	UNP Q96C86
D	1	GLY	-	EXPRESSION TAG	UNP Q96C86

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is 2-[(2,4-diaminoquinazolin-5-yl)oxy]methyl} benzenesulfonic acid (three-letter code: 30S) (formula: C₁₅H₁₄N₄O₄S).

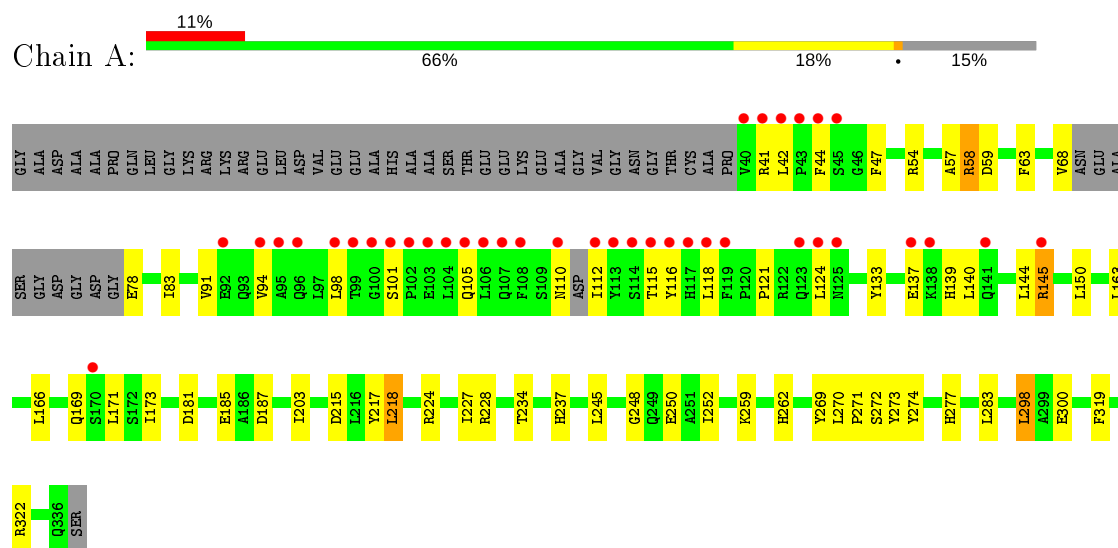


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	S	13	0
			36	15	13	4	3	1		
3	C	1	Total	C	N	O	S		0	0
			23	15	4	3	1			

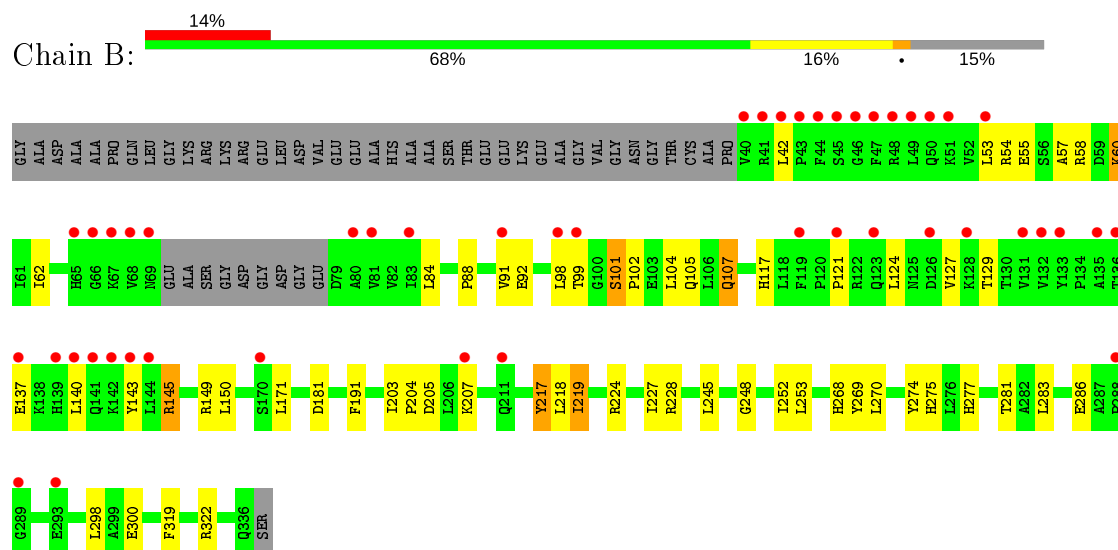
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: m7GpppX diphosphatase

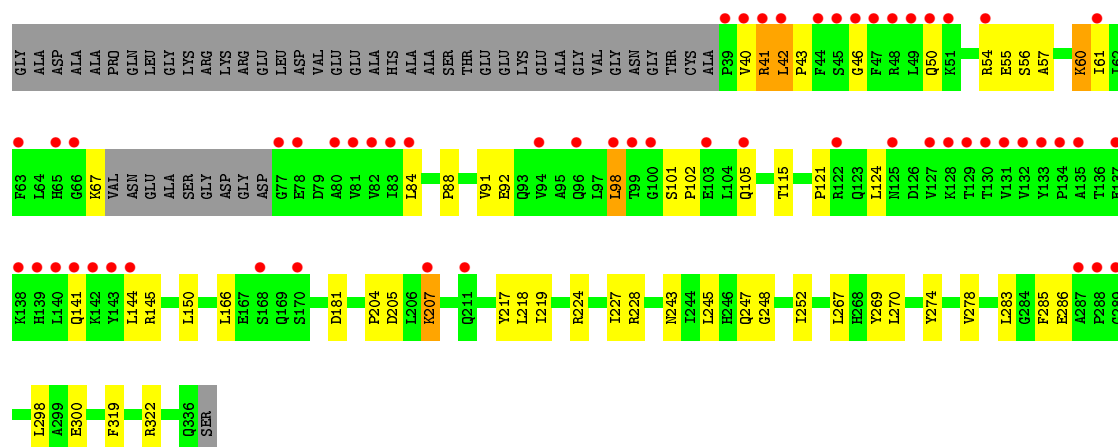


• Molecule 1: m7GpppX diphosphatase

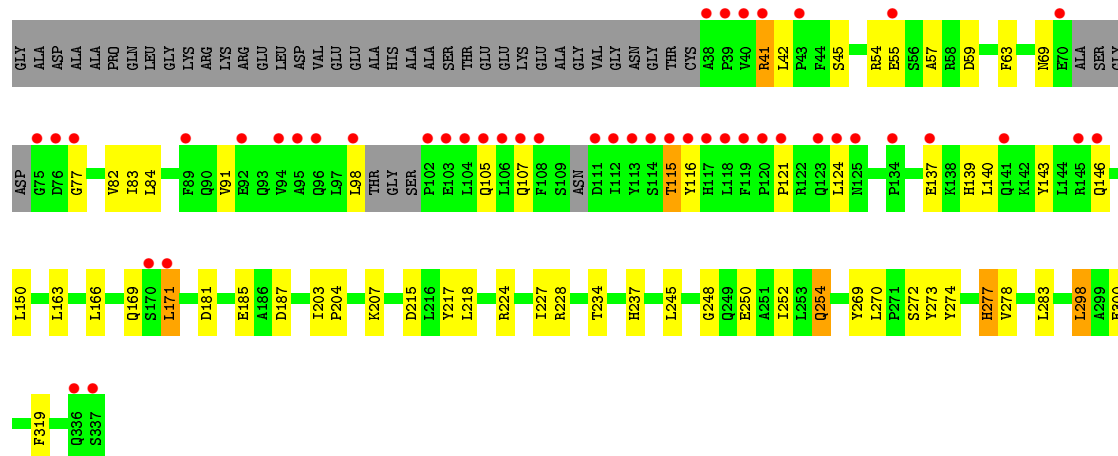


• Molecule 1: m7GpppX diphosphatase





• Molecule 1: m7GpppX diphosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.37Å 105.61Å 137.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.68 – 2.90 38.67 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.68-2.90) 99.9 (38.67-2.90)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.90Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.209 , 0.251 0.235 , 0.288	Depositor DCC
R_{free} test set	1683 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 68.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9655	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.90 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.9842e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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, 3OS

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	0.51	0/2432	0.74	2/3298 (0.1%)
1	B	0.49	0/2440	0.74	2/3311 (0.1%)
1	C	0.50	0/2446	0.75	3/3318 (0.1%)
1	D	0.51	0/2466	0.75	1/3342 (0.0%)
All	All	0.50	0/9784	0.74	8/13269 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	269	TYR	C-N-CA	8.02	141.76	121.70
1	A	269	TYR	C-N-CA	7.83	141.28	121.70
1	B	269	TYR	C-N-CA	7.78	141.14	121.70
1	C	269	TYR	C-N-CA	7.65	140.82	121.70
1	C	286	GLU	C-N-CA	5.26	134.85	121.70
1	B	269	TYR	N-CA-CB	5.17	119.90	110.60
1	A	269	TYR	N-CA-CB	5.08	119.75	110.60
1	C	269	TYR	N-CA-CB	5.03	119.65	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2377	0	2383	38	0
1	B	2384	0	2388	40	0
1	C	2389	0	2390	27	0
1	D	2411	0	2405	37	0
2	A	5	0	0	1	0
2	B	10	0	0	3	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	B	23	13	13	0	0
3	C	23	0	13	2	0
All	All	9642	13	9592	113	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 6.

All (113) 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:B:219:ILE:HD11	1:B:277:HIS:HB2	1.24	1.09
1:B:54:ARG:HD3	1:B:145:ARG:HE	1.32	0.91
1:B:275:HIS:O	1:B:277:HIS:HD2	1.63	0.80
1:A:234:THR:H	1:A:237:HIS:HD2	1.33	0.75
1:D:234:THR:H	1:D:237:HIS:HD2	1.35	0.73
1:C:205:ASP:OD1	1:C:207:LYS:HG2	1.87	0.73
1:D:273:TYR:HB2	1:D:277:HIS:HE1	1.58	0.69
1:B:277:HIS:HE1	2:B:402:PO4:O4	1.76	0.68
1:D:41:ARG:HD2	1:D:42:LEU:H	1.61	0.66
1:C:150:LEU:HD13	1:D:283:LEU:HD11	1.77	0.65
1:B:205:ASP:OD1	1:B:207:LYS:HG2	1.97	0.65
1:A:42:LEU:HG	1:B:104:LEU:CD1	2.27	0.64
1:B:191:PHE:CZ	1:D:254:GLN:HB2	2.34	0.63
1:B:219:ILE:HD11	1:B:277:HIS:CB	2.17	0.62
1:C:102:PRO:HD2	1:D:41:ARG:HD3	1.81	0.62
1:C:60:LYS:HE2	1:D:59:ASP:O	2.01	0.61
1:B:219:ILE:CD1	1:B:277:HIS:HB2	2.16	0.61
1:C:57:ALA:HB1	1:D:215:ASP:HB3	1.82	0.60
1:A:42:LEU:HG	1:B:104:LEU:HD13	1.83	0.59
1:A:228:ARG:HB3	1:A:274:TYR:HB3	1.83	0.59
1:C:41:ARG:HB3	1:D:116:TYR:HE2	1.68	0.59
1:B:228:ARG:HB3	1:B:274:TYR:HB3	1.85	0.59
1:A:94:VAL:HG11	1:B:62:ILE:HD11	1.85	0.59
1:C:228:ARG:HB3	1:C:274:TYR:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:GLN:HE21	1:C:144:LEU:HD22	1.69	0.58
1:D:273:TYR:HB2	1:D:277:HIS:CE1	2.38	0.56
1:C:204:PRO:HA	1:C:218:LEU:HG	1.87	0.56
1:C:98:LEU:HD11	1:D:84:LEU:HD11	1.88	0.55
1:D:207:LYS:HD2	1:D:217:TYR:HE2	1.71	0.55
1:B:204:PRO:HA	1:B:218:LEU:HG	1.88	0.55
1:A:215:ASP:HB3	1:B:57:ALA:HB1	1.88	0.55
1:B:224:ARG:HD3	1:B:227:ILE:HD11	1.89	0.54
1:A:283:LEU:HD11	1:B:150:LEU:HD13	1.89	0.53
1:C:217:TYR:HD2	1:C:285:PHE:HE2	1.57	0.53
1:A:63:PHE:CD2	1:A:83:ILE:HD12	2.44	0.52
1:A:215:ASP:HB3	1:B:57:ALA:CB	2.39	0.52
1:C:224:ARG:HD3	1:C:227:ILE:HD11	1.92	0.52
1:D:63:PHE:CD2	1:D:83:ILE:HD12	2.45	0.51
1:A:59:ASP:O	1:B:60:LYS:HE2	2.11	0.51
1:C:42:LEU:HD13	1:C:42:LEU:H	1.74	0.51
1:C:98:LEU:CD1	1:D:84:LEU:HD11	2.41	0.51
1:A:44:PHE:HB2	1:A:68:VAL:HB	1.93	0.50
1:A:54:ARG:HH11	1:A:145:ARG:HG2	1.76	0.50
1:D:203:ILE:HG13	1:D:204:PRO:HD2	1.93	0.50
1:A:224:ARG:HD3	1:A:227:ILE:HD11	1.93	0.50
1:C:283:LEU:HD11	1:D:150:LEU:HD13	1.92	0.50
1:D:224:ARG:HD3	1:D:227:ILE:HD11	1.93	0.49
3:C:401:3OS:H3	1:D:143:TYR:OH	2.13	0.49
1:A:185:GLU:HB3	1:A:203:ILE:HD11	1.93	0.49
1:C:57:ALA:CB	1:D:215:ASP:HB3	2.42	0.48
1:C:56:SER:HB3	1:C:61:ILE:HG22	1.94	0.48
1:A:145:ARG:HA	1:A:145:ARG:HE	1.79	0.48
1:D:107:GLN:HB3	1:D:115:THR:HB	1.96	0.48
1:D:228:ARG:HB3	1:D:274:TYR:HB3	1.95	0.47
1:B:54:ARG:HB2	1:B:145:ARG:HB2	1.96	0.47
1:B:107:GLN:HG3	1:B:117:HIS:NE2	2.30	0.47
1:B:121:PRO:HD2	1:B:124:LEU:HD13	1.97	0.47
1:D:121:PRO:HD2	1:D:124:LEU:HD13	1.97	0.47
1:B:42:LEU:HD22	1:B:42:LEU:H	1.80	0.46
1:B:277:HIS:CE1	2:B:402:PO4:O4	2.64	0.46
1:B:268:HIS:HB2	2:B:402:PO4:O2	2.15	0.46
1:A:248:GLY:O	1:A:252:ILE:HG12	2.16	0.46
1:C:248:GLY:O	1:C:252:ILE:HG12	2.16	0.46
1:D:185:GLU:HB3	1:D:203:ILE:HD11	1.98	0.46
1:D:245:LEU:HD13	1:D:298:LEU:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:TYR:HB2	1:B:129:THR:HB	1.97	0.45
1:D:248:GLY:O	1:D:252:ILE:HG12	2.17	0.45
1:C:121:PRO:HD2	1:C:124:LEU:HD13	1.98	0.45
1:D:41:ARG:HH22	1:D:45:SER:HA	1.82	0.45
1:B:270:LEU:HD22	1:B:319:PHE:CE2	2.51	0.45
1:A:245:LEU:HD13	1:A:298:LEU:HB3	1.97	0.45
1:B:248:GLY:O	1:B:252:ILE:HG12	2.17	0.45
1:A:121:PRO:HD2	1:A:124:LEU:HD13	1.98	0.45
1:D:270:LEU:HD22	1:D:319:PHE:CE2	2.51	0.45
1:A:137:GLU:HA	1:A:140:LEU:HD12	1.98	0.44
1:A:150:LEU:HD13	1:B:283:LEU:HD11	1.98	0.44
1:C:54:ARG:NE	1:C:145:ARG:HD3	2.33	0.44
1:D:54:ARG:HB3	1:D:63:PHE:CD1	2.53	0.44
1:A:270:LEU:HD22	1:A:319:PHE:CE2	2.52	0.44
1:C:270:LEU:HD22	1:C:319:PHE:CE2	2.52	0.44
1:A:118:LEU:HB3	1:B:127:VAL:HB	2.00	0.43
1:B:137:GLU:HA	1:B:140:LEU:HD12	2.00	0.43
1:A:271:PRO:HA	2:A:401:PO4:O2	2.19	0.43
1:A:57:ALA:HA	1:B:88:PRO:HB3	2.01	0.43
1:D:54:ARG:HB2	1:D:143:TYR:O	2.18	0.43
1:C:50:GLN:HE21	1:C:67:LYS:HE3	1.83	0.43
1:A:41:ARG:HG3	1:A:42:LEU:N	2.34	0.43
1:A:47:PHE:O	1:B:99:THR:HG22	2.18	0.43
1:D:270:LEU:HA	1:D:270:LEU:HD12	1.83	0.43
1:C:150:LEU:HD13	1:D:283:LEU:CD1	2.47	0.43
1:A:218:LEU:HD13	1:A:252:ILE:CD1	2.48	0.43
1:A:173:ILE:HD12	1:A:273:TYR:HB3	2.01	0.42
1:C:267:LEU:HG	1:C:278:VAL:HG22	2.01	0.42
1:D:137:GLU:HA	1:D:140:LEU:HD12	2.01	0.42
1:B:92:GLU:HG2	1:B:92:GLU:H	1.55	0.42
1:A:54:ARG:H	1:A:63:PHE:HB2	1.84	0.41
1:A:78:GLU:HB3	1:A:133:TYR:HE1	1.85	0.41
1:D:278:VAL:HG11	1:D:298:LEU:HD12	2.03	0.41
1:A:270:LEU:HD12	1:A:270:LEU:HA	1.84	0.41
1:A:42:LEU:HG	1:B:104:LEU:HD11	1.99	0.41
1:C:243:ASN:O	1:C:247:GLN:HB2	2.20	0.41
1:B:53:LEU:HD22	1:B:143:TYR:HB2	2.02	0.41
1:A:218:LEU:HD13	1:A:252:ILE:HD12	2.02	0.41
1:A:41:ARG:HH22	1:A:42:LEU:HD12	1.86	0.41
1:B:101:SER:HA	1:B:102:PRO:HD3	1.98	0.41
1:C:88:PRO:HB3	1:D:57:ALA:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:HIS:NE2	1:B:149:ARG:CZ	2.84	0.41
1:B:217:TYR:CD1	1:B:281:THR:HG22	2.56	0.40
1:C:98:LEU:HD11	1:D:82:VAL:HG11	2.03	0.40
1:A:94:VAL:HG11	1:B:62:ILE:CD1	2.51	0.40
1:A:58:ARG:HH22	1:B:286:GLU:HB2	1.85	0.40
3:C:401:30S:N2	1:D:185:GLU:OE2	2.55	0.40
1:D:63:PHE:HD2	1:D:83:ILE:HD12	1.86	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	281/337 (83%)	270 (96%)	8 (3%)	3 (1%)	14	42
1	B	284/337 (84%)	271 (95%)	12 (4%)	1 (0%)	34	66
1	C	285/337 (85%)	264 (93%)	17 (6%)	4 (1%)	11	36
1	D	284/337 (84%)	267 (94%)	13 (5%)	4 (1%)	11	36
All	All	1134/1348 (84%)	1072 (94%)	50 (4%)	12 (1%)	14	42

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	GLN
1	D	169	GLN
1	B	91	VAL
1	C	43	PRO
1	C	46	GLY
1	A	91	VAL
1	A	171	LEU
1	C	91	VAL

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Mol	Chain	Res	Type
1	C	40	VAL
1	D	91	VAL
1	D	171	LEU
1	D	77	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	261/294 (89%)	238 (91%)	23 (9%)	10	30
1	B	262/294 (89%)	243 (93%)	19 (7%)	14	38
1	C	262/294 (89%)	244 (93%)	18 (7%)	15	41
1	D	264/294 (90%)	244 (92%)	20 (8%)	13	36
All	All	1049/1176 (89%)	969 (92%)	80 (8%)	13	36

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ARG
1	A	98	LEU
1	A	101	SER
1	A	105	GLN
1	A	110	ASN
1	A	112	ILE
1	A	115	THR
1	A	139	HIS
1	A	144	LEU
1	A	145	ARG
1	A	163	LEU
1	A	166	LEU
1	A	181	ASP
1	A	187	ASP
1	A	217	TYR
1	A	218	LEU
1	A	250	GLU

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Mol	Chain	Res	Type
1	A	259	LYS
1	A	272	SER
1	A	277	HIS
1	A	298	LEU
1	A	300	GLU
1	A	322	ARG
1	B	55	GLU
1	B	58	ARG
1	B	60	LYS
1	B	84	LEU
1	B	98	LEU
1	B	101	SER
1	B	105	GLN
1	B	107	GLN
1	B	145	ARG
1	B	171	LEU
1	B	181	ASP
1	B	203	ILE
1	B	217	TYR
1	B	219	ILE
1	B	245	LEU
1	B	253	LEU
1	B	298	LEU
1	B	300	GLU
1	B	322	ARG
1	C	41	ARG
1	C	42	LEU
1	C	55	GLU
1	C	60	LYS
1	C	84	LEU
1	C	92	GLU
1	C	98	LEU
1	C	101	SER
1	C	105	GLN
1	C	115	THR
1	C	166	LEU
1	C	181	ASP
1	C	207	LYS
1	C	219	ILE
1	C	245	LEU
1	C	298	LEU
1	C	300	GLU

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Mol	Chain	Res	Type
1	C	322	ARG
1	D	41	ARG
1	D	55	GLU
1	D	69	ASN
1	D	98	LEU
1	D	105	GLN
1	D	115	THR
1	D	139	HIS
1	D	146	GLN
1	D	163	LEU
1	D	166	LEU
1	D	171	LEU
1	D	181	ASP
1	D	187	ASP
1	D	218	LEU
1	D	250	GLU
1	D	254	GLN
1	D	272	SER
1	D	277	HIS
1	D	298	LEU
1	D	300	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	146	GLN
1	A	237	HIS
1	B	169	GLN
1	B	209	ASN
1	B	211	GLN
1	B	212	GLN
1	B	277	HIS
1	C	50	GLN
1	C	141	GLN
1	C	209	ASN
1	C	211	GLN
1	C	212	GLN
1	D	237	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

9 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$
2	PO4	A	401	-	4,4,4	2.45	2 (50%)	6,6,6	0.97	0
2	PO4	C	402	-	4,4,4	1.82	0	6,6,6	0.84	0
3	30S	B	401	1	22,25,26	0.51	0	31,35,38	1.20	3 (9%)
2	PO4	D	401	-	4,4,4	2.49	2 (50%)	6,6,6	0.67	0
2	PO4	B	402	-	4,4,4	2.42	1 (25%)	6,6,6	0.77	0
2	PO4	D	402	-	4,4,4	1.95	1 (25%)	6,6,6	0.75	0
2	PO4	C	403	-	4,4,4	2.56	2 (50%)	6,6,6	0.45	0
3	30S	C	401	1	22,25,26	0.56	1 (4%)	31,35,38	1.21	3 (9%)
2	PO4	B	403	-	4,4,4	2.51	2 (50%)	6,6,6	1.32	1 (16%)

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	30S	C	401	1	-	1/9/9/11	0/3/3/3
3	30S	B	401	1	-	1/9/9/11	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	403	PO4	P-O1	4.25	1.60	1.50
2	B	403	PO4	P-O1	4.12	1.60	1.50
2	D	401	PO4	P-O1	4.10	1.60	1.50
2	B	402	PO4	P-O1	4.03	1.60	1.50
2	A	401	PO4	P-O1	4.00	1.60	1.50
2	B	403	PO4	P-O4	2.05	1.60	1.54
2	D	401	PO4	P-O3	2.04	1.60	1.54
2	D	402	PO4	P-O1	2.04	1.55	1.50
2	C	403	PO4	P-O3	2.01	1.60	1.54
3	C	401	30S	C5-C4	2.00	1.41	1.40
2	A	401	PO4	P-O3	2.00	1.60	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	401	30S	C8-C13-C14	4.72	129.75	127.36
3	B	401	30S	C8-C13-C14	4.64	129.71	127.36
3	C	401	30S	C14-C13-C12	-2.65	113.05	114.86
3	B	401	30S	C14-C13-C12	-2.48	113.17	114.86
2	B	403	PO4	O4-P-O2	2.35	115.52	107.97
3	B	401	30S	C5-C4-S1	2.31	127.46	121.18
3	C	401	30S	C5-C4-S1	2.31	127.45	121.18

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	401	30S	C5-C4-S1-O3
3	C	401	30S	C5-C4-S1-O3

There are no ring outliers.

3 monomers are involved in 6 short contacts:

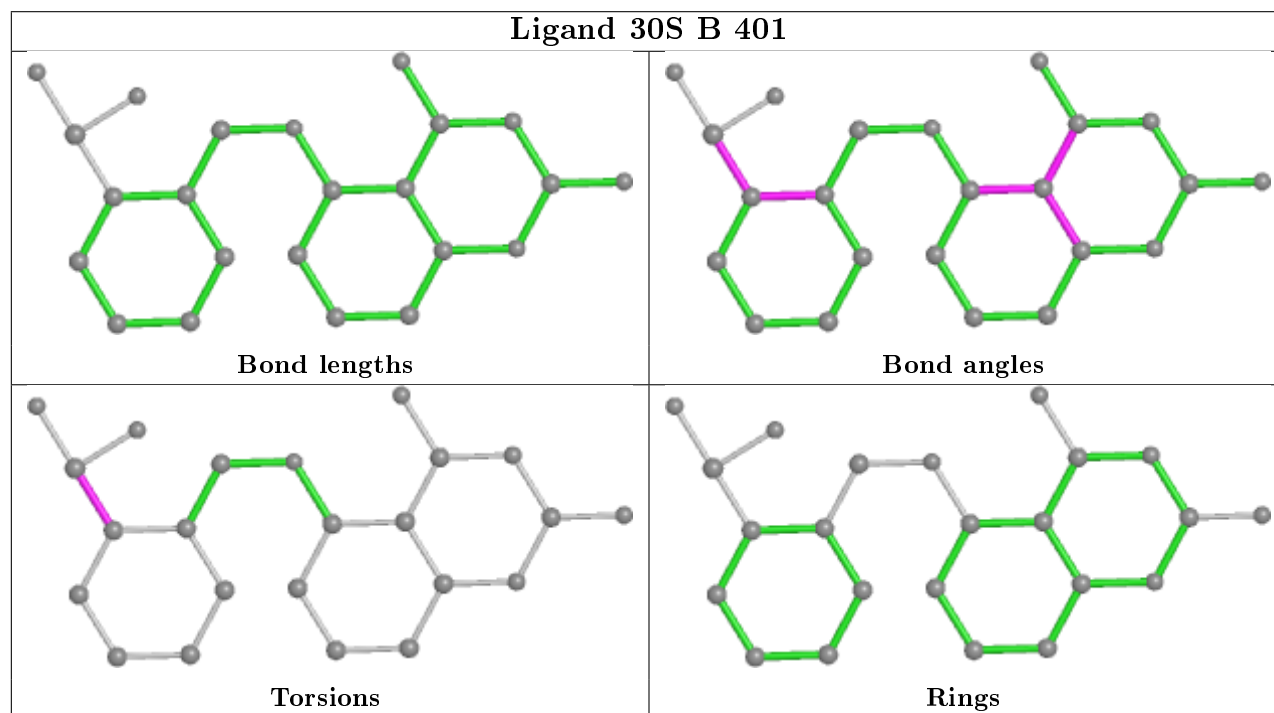
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	PO4	1	0
2	B	402	PO4	3	0

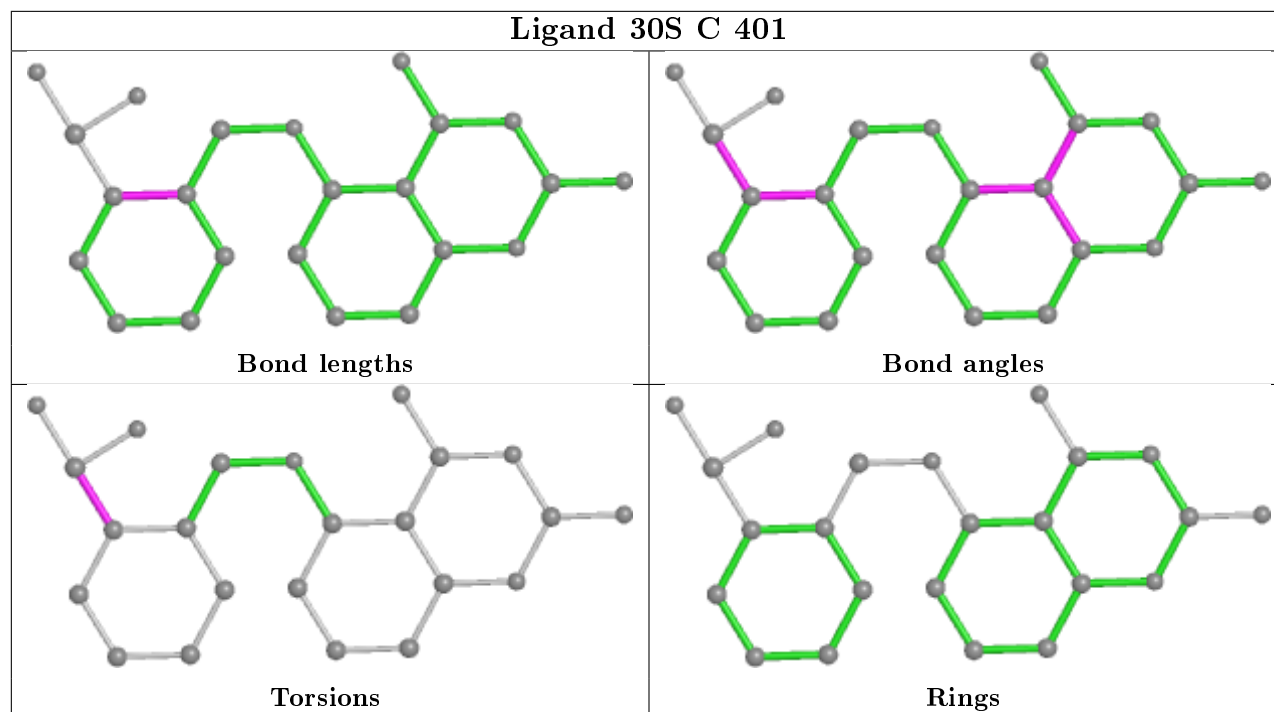
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	401	30S	2	0

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.





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	287/337 (85%)	0.67	38 (13%) 3 2	21, 52, 140, 169	0
1	B	288/337 (85%)	0.75	47 (16%) 1 1	23, 53, 147, 208	0
1	C	289/337 (85%)	0.73	57 (19%) 1 0	21, 52, 147, 188	0
1	D	292/337 (86%)	0.67	46 (15%) 2 1	20, 50, 144, 180	0
All	All	1156/1348 (85%)	0.70	188 (16%) 1 1	20, 52, 147, 208	0

All (188) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	43	PRO	10.5
1	B	47	PHE	10.1
1	D	111	ASP	9.9
1	D	170	SER	9.6
1	A	113	TYR	9.4
1	D	39	PRO	8.8
1	D	113	TYR	8.5
1	A	112	ILE	8.4
1	B	44	PHE	8.2
1	C	40	VAL	8.2
1	B	45	SER	7.7
1	C	39	PRO	7.6
1	A	102	PRO	7.5
1	A	101	SER	7.2
1	C	41	ARG	7.0
1	A	107	GLN	7.0
1	D	118	LEU	6.9
1	D	95	ALA	6.9
1	D	38	ALA	6.8
1	B	40	VAL	6.7
1	A	104	LEU	6.7

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Mol	Chain	Res	Type	RSRZ
1	D	337	SER	6.7
1	B	42	LEU	6.5
1	B	69	ASN	6.4
1	D	94	VAL	6.2
1	A	41	ARG	6.1
1	A	103	GLU	6.1
1	B	68	VAL	5.9
1	C	129	THR	5.9
1	D	112	ILE	5.8
1	A	98	LEU	5.8
1	A	110	ASN	5.7
1	A	99	THR	5.4
1	B	142	LYS	5.4
1	B	132	VAL	5.4
1	A	43	PRO	5.3
1	B	66	GLY	5.3
1	C	42	LEU	5.2
1	A	40	VAL	5.2
1	A	106	LEU	5.2
1	C	47	PHE	5.1
1	C	65	HIS	5.1
1	B	50	GLN	5.1
1	B	139	HIS	5.1
1	A	94	VAL	5.1
1	C	66	GLY	5.0
1	D	104	LEU	4.9
1	A	123	GLN	4.7
1	B	80	ALA	4.7
1	A	170	SER	4.7
1	D	41	ARG	4.5
1	B	143	TYR	4.5
1	B	135	ALA	4.5
1	B	51	LYS	4.5
1	D	107	GLN	4.5
1	B	41	ARG	4.5
1	B	133	TYR	4.5
1	C	77	GLY	4.4
1	D	124	LEU	4.3
1	D	108	PHE	4.3
1	D	114	SER	4.3
1	C	130	THR	4.2
1	A	116	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	138	LYS	4.2
1	D	121	PRO	4.1
1	A	100	GLY	4.0
1	D	120	PRO	3.9
1	A	115	THR	3.9
1	D	102	PRO	3.9
1	C	49	LEU	3.9
1	D	116	TYR	3.9
1	D	75	GLY	3.8
1	C	141	GLN	3.8
1	B	48	ARG	3.7
1	A	119	PHE	3.7
1	B	131	VAL	3.7
1	C	48	ARG	3.7
1	C	288	PRO	3.6
1	C	170	SER	3.6
1	C	50	GLN	3.6
1	B	46	GLY	3.6
1	C	287	ALA	3.6
1	C	144	LEU	3.6
1	C	44	PHE	3.6
1	A	118	LEU	3.5
1	C	131	VAL	3.5
1	D	125	ASN	3.5
1	C	81	VAL	3.5
1	A	114	SER	3.5
1	A	108	PHE	3.5
1	C	137	GLU	3.4
1	B	49	LEU	3.4
1	D	98	LEU	3.4
1	D	115	THR	3.4
1	D	106	LEU	3.3
1	B	81	VAL	3.3
1	B	170	SER	3.2
1	C	45	SER	3.2
1	C	122	ARG	3.2
1	C	134	PRO	3.2
1	C	139	HIS	3.2
1	A	141	GLN	3.2
1	B	123	GLN	3.2
1	D	119	PHE	3.2
1	C	103	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	65	HIS	3.2
1	C	211	GLN	3.2
1	D	134	PRO	3.2
1	A	125	ASN	3.1
1	D	76	ASP	3.1
1	C	46	GLY	3.1
1	B	67	LYS	3.1
1	D	40	VAL	3.1
1	C	133	TYR	3.0
1	B	141	GLN	3.0
1	A	145	ARG	2.9
1	B	126	ASP	2.9
1	C	143	TYR	2.9
1	B	211	GLN	2.9
1	D	89	PHE	2.9
1	C	168	SER	2.9
1	A	45	SER	2.8
1	C	140	LEU	2.8
1	A	124	LEU	2.8
1	D	96	GLN	2.8
1	B	288	PRO	2.7
1	A	42	LEU	2.7
1	C	105	GLN	2.7
1	B	207	LYS	2.6
1	D	105	GLN	2.6
1	B	99	THR	2.6
1	C	127	VAL	2.6
1	C	54	ARG	2.6
1	D	145	ARG	2.6
1	C	63	PHE	2.6
1	D	123	GLN	2.5
1	B	137	GLU	2.5
1	A	137	GLU	2.5
1	B	136	THR	2.5
1	D	141	GLN	2.5
1	D	117	HIS	2.5
1	C	128	LYS	2.5
1	C	207	LYS	2.5
1	B	83	ILE	2.5
1	B	119	PHE	2.5
1	C	80	ALA	2.5
1	B	144	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	103	GLU	2.5
1	D	336	GLN	2.5
1	B	140	LEU	2.5
1	C	51	LYS	2.5
1	B	121	PRO	2.4
1	A	92	GLU	2.4
1	C	98	LEU	2.4
1	C	83	ILE	2.4
1	C	96	GLN	2.4
1	D	55	GLU	2.3
1	D	137	GLU	2.3
1	C	142	LYS	2.3
1	C	61	ILE	2.3
1	B	91	VAL	2.3
1	D	43	PRO	2.3
1	A	105	GLN	2.3
1	C	94	VAL	2.3
1	D	70	GLU	2.3
1	A	96	GLN	2.3
1	B	53	LEU	2.2
1	D	77	GLY	2.2
1	D	171	LEU	2.2
1	C	135	ALA	2.2
1	C	125	ASN	2.2
1	D	146	GLN	2.1
1	C	82	VAL	2.1
1	B	98	LEU	2.1
1	C	78	GLU	2.1
1	C	132	VAL	2.1
1	A	138	LYS	2.1
1	B	289	GLY	2.1
1	A	44	PHE	2.1
1	B	128	LYS	2.1
1	C	100	GLY	2.1
1	C	84	LEU	2.0
1	A	95	ALA	2.0
1	D	92	GLU	2.0
1	B	293	GLU	2.0
1	A	117	HIS	2.0
1	C	289	GLY	2.0
1	C	99	THR	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

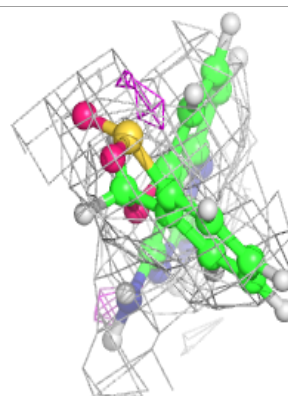
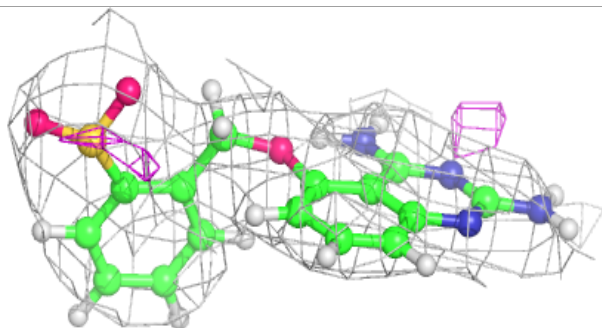
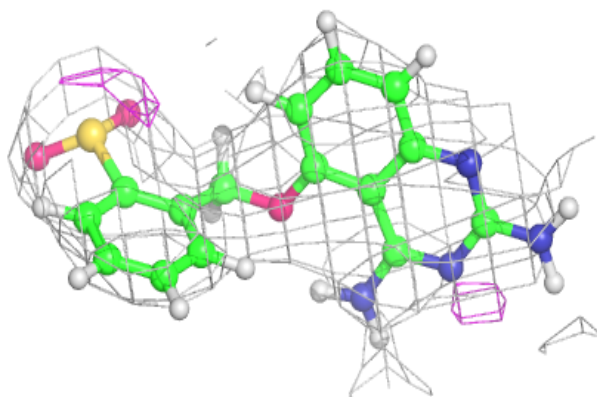
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PO4	B	403	5/5	0.83	0.39	67,67,70,71	0
2	PO4	C	403	5/5	0.89	0.32	64,65,70,72	0
3	30S	B	401	23/24	0.93	0.16	47,51,61,63	13
2	PO4	D	401	5/5	0.93	0.24	67,68,70,74	0
2	PO4	A	401	5/5	0.94	0.29	57,59,60,61	0
2	PO4	D	402	5/5	0.94	0.34	71,74,76,78	0
3	30S	C	401	23/24	0.95	0.20	39,45,62,65	0
2	PO4	B	402	5/5	0.97	0.14	41,42,42,46	0
2	PO4	C	402	5/5	0.98	0.18	39,43,44,44	0

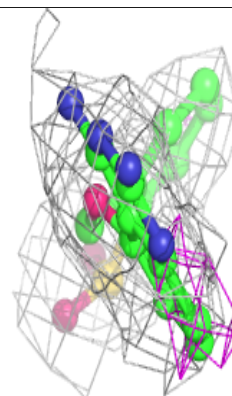
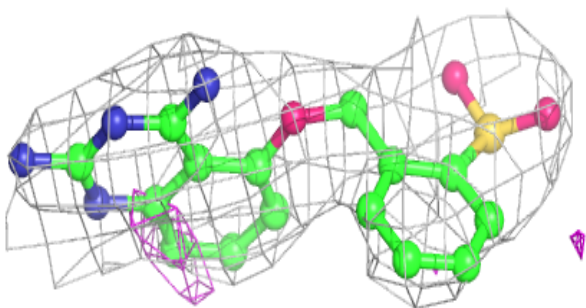
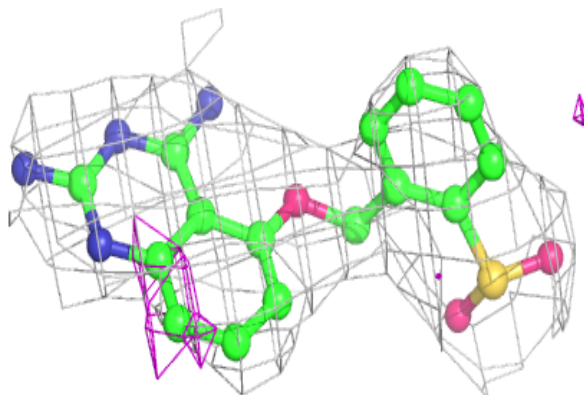
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 30S 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 30S 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.