



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:37 PM UTC

PDB ID : 6YGJ / pdb_00006ygj
Title : small-molecule stabilizer of 14-3-3 and the Carbohydrate Response Element Binding Protein (ChREBP) protein-protein interaction
Authors : Ottmann, C.; Visser, E.J.
Deposited on : 2020-03-27
Resolution : 2.07 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

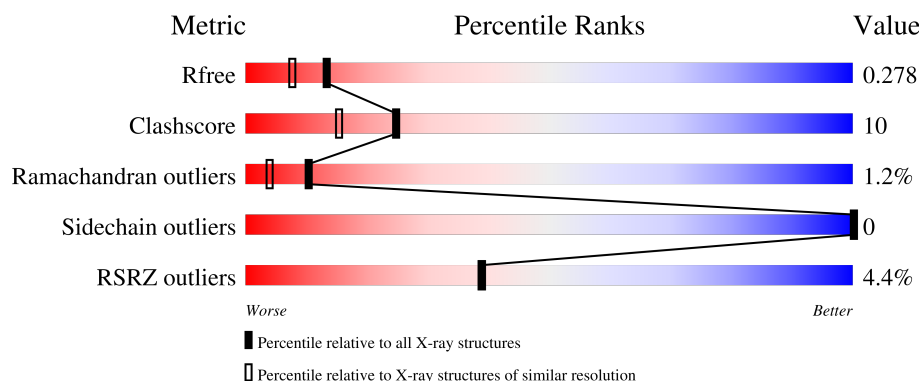
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	229	3% 82% 17% ..
2	B	21	62% 19% 14% 5%
2	I	21	5% 67% 29% 5%
3	F	231	6% 72% 26% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8135 atoms, of which 4021 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 14-3-3 protein beta/alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	227	Total	C	H	N	O	S	0	0	0
			3633	1145	1806	306	368	8			

- Molecule 2 is a protein called Carbohydrate-responsive element-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	20	Total	C	H	N	O	0	0	0
			369	121	184	36	28			
2	I	20	Total	C	H	N	O	0	0	0
			369	121	184	36	28			

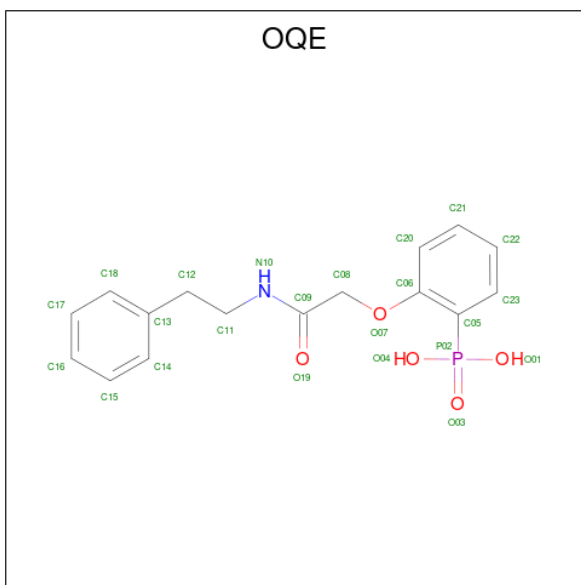
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	136	GLN	LYS	conflict	UNP Q9NP71
I	136	GLN	LYS	conflict	UNP Q9NP71

- Molecule 3 is a protein called 14-3-3 protein beta/alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	F	228	Total	C	H	N	O	S	0	0	0
			3650	1150	1815	307	369	9			

- Molecule 4 is [2-[2-oxidanylidene-2-(2-phenylethylamino)ethoxy]phenyl]phosphonic acid (CCD ID: OQE) (formula: C₁₆H₁₈NO₅P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	
			39	16	16	1	5	1	0
4	F	1	Total	C	H	N	O	P	
			39	16	16	1	5	1	0

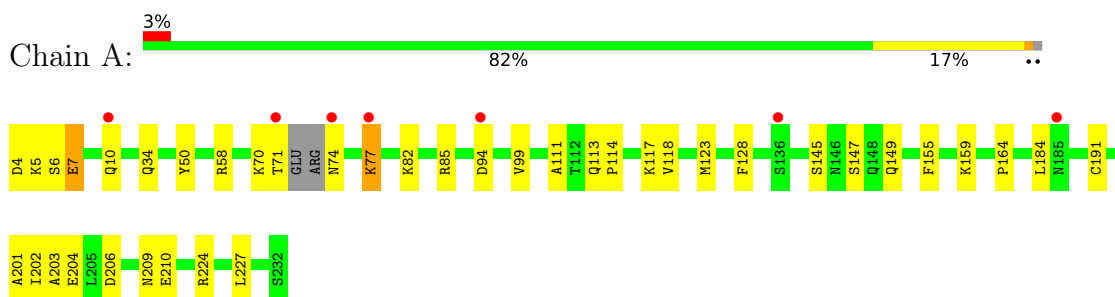
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	12	Total	O		
			12	12	0	0
5	B	4	Total	O		
			4	4	0	0
5	F	19	Total	O		
			19	19	0	0
5	I	1	Total	O		
			1	1	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.

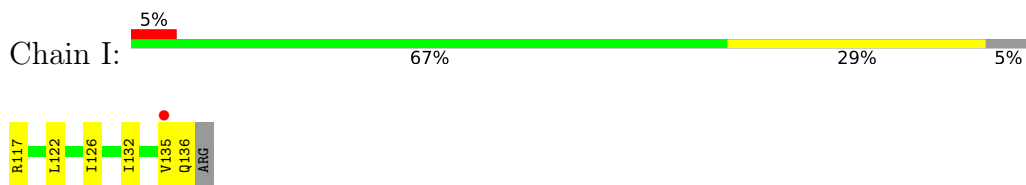
- Molecule 1: 14-3-3 protein beta/alpha



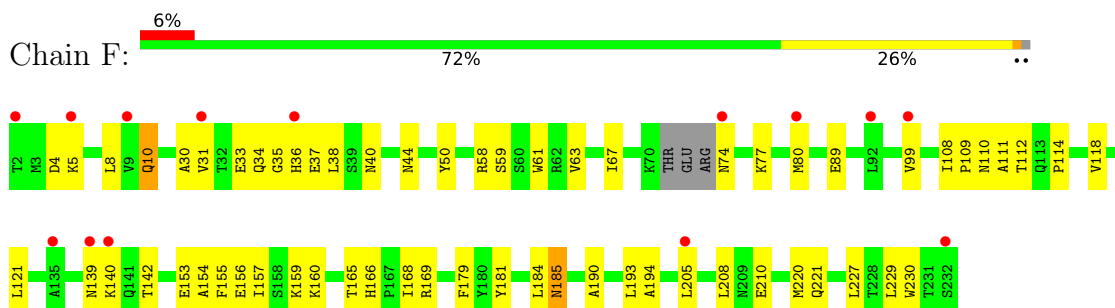
- Molecule 2: Carbohydrate-responsive element-binding protein



- Molecule 2: Carbohydrate-responsive element-binding protein



- Molecule 3: 14-3-3 protein beta/alpha



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.79Å 76.69Å 90.29Å 90.00° 119.22° 90.00°	Depositor
Resolution (Å)	57.30 – 2.07 57.30 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.9 (57.30-2.07) 99.9 (57.30-2.07)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.07Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.221 , 0.277 0.226 , 0.278	Depositor DCC
R_{free} test set	1778 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	57.1	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 54.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8135	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.76% 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: OQE

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	0.71	4/1853 (0.2%)	0.88	5/2494 (0.2%)
2	B	0.59	0/190	0.88	0/257
2	I	0.58	0/190	0.65	0/257
3	F	0.75	3/1861 (0.2%)	0.85	5/2504 (0.2%)
All	All	0.72	7/4094 (0.2%)	0.86	10/5512 (0.2%)

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
2	B	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	LYS	CB-CG	-8.30	1.27	1.52
3	F	10	GLN	CD-NE2	-8.19	1.16	1.33
1	A	7	GLU	CD-OE1	7.21	1.39	1.25
3	F	220	MET	SD-CE	-5.62	1.65	1.79
1	A	94	ASP	CB-CG	5.39	1.65	1.52
3	F	80	MET	CG-SD	-5.26	1.67	1.80
1	A	123	MET	SD-CE	-5.05	1.67	1.79

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	LYS	CD-CE-NZ	-16.41	59.40	111.90
3	F	10	GLN	CA-CB-CG	-9.12	95.86	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	10	GLN	OE1-CD-NE2	-8.79	113.81	122.60
3	F	80	MET	CA-CB-CG	-7.09	99.92	114.10
3	F	10	GLN	CG-CD-NE2	6.50	126.15	116.40
1	A	77	LYS	CB-CG-CD	-6.38	96.62	111.30
3	F	10	GLN	CB-CG-CD	5.62	122.15	112.60
1	A	7	GLU	CB-CG-CD	-5.46	103.32	112.60
1	A	7	GLU	CA-CB-CG	-5.39	103.31	114.10
1	A	7	GLU	CG-CD-OE1	5.37	130.75	118.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	117	ARG	Peptide
2	B	118	ASP	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	1827	1806	1804	27	3
2	B	185	184	184	6	0
2	I	185	184	184	7	0
3	F	1835	1815	1815	42	3
4	A	23	16	0	2	0
4	F	23	16	0	2	0
5	A	12	0	0	0	0
5	B	4	0	0	0	0
5	F	19	0	0	0	0
5	I	1	0	0	0	0
All	All	4114	4021	3987	79	3

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 10.

All (79) 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:5:LYS:C	1:A:7:GLU:H	1.98	0.71
1:A:7:GLU:HA	1:A:7:GLU:OE1	1.93	0.68
3:F:165:THR:HB	3:F:208:LEU:HD13	1.75	0.68
3:F:139:ASN:O	3:F:142:THR:N	2.29	0.66
2:B:117:ARG:HH12	2:B:119:LYS:HE2	1.62	0.63
3:F:205:LEU:O	3:F:205:LEU:HG	2.00	0.61
1:A:5:LYS:O	1:A:7:GLU:N	2.34	0.60
3:F:34:GLN:O	3:F:36:HIS:ND1	2.32	0.60
1:A:145:SER:O	1:A:149:GLN:HG2	2.03	0.59
1:A:58:ARG:HH12	4:A:301:OQE:C05	2.16	0.58
1:A:4:ASP:O	1:A:7:GLU:HB2	2.05	0.57
2:B:117:ARG:CZ	2:B:117:ARG:HB3	2.32	0.57
1:A:5:LYS:C	1:A:7:GLU:N	2.62	0.56
3:F:10:GLN:CD	3:F:10:GLN:O	2.48	0.56
3:F:108:ILE:HB	3:F:109:PRO:HD3	1.87	0.55
1:A:71:THR:HG22	1:A:71:THR:O	2.06	0.55
3:F:156:GLU:O	3:F:160:LYS:HE2	2.06	0.55
3:F:194:ALA:HB3	3:F:227:LEU:HD21	1.88	0.54
3:F:10:GLN:O	3:F:10:GLN:OE1	2.24	0.54
3:F:40:ASN:O	3:F:44:ASN:OD1	2.26	0.53
3:F:37:GLU:HG3	3:F:38:LEU:H	1.74	0.53
3:F:229:LEU:HD21	2:I:132:ILE:HD12	1.90	0.52
3:F:37:GLU:HG3	3:F:38:LEU:N	2.24	0.52
1:A:58:ARG:NH1	4:A:301:OQE:C23	2.73	0.52
1:A:164:PRO:O	1:A:201:ALA:HB2	2.09	0.52
2:B:124:ASN:O	2:B:128:ARG:HG2	2.10	0.51
1:A:74:ASN:OD1	1:A:77:LYS:CB	2.59	0.51
3:F:153:GLU:O	3:F:157:ILE:HG13	2.10	0.51
3:F:155:PHE:CE1	3:F:159:LYS:HE3	2.45	0.50
3:F:221:GLN:NE2	2:I:117:ARG:HH21	2.09	0.50
1:A:10:GLN:O	1:A:10:GLN:HG2	2.12	0.50
3:F:229:LEU:HD21	2:I:132:ILE:CD1	2.42	0.49
3:F:59:SER:O	3:F:63:VAL:HG23	2.13	0.49
1:A:184:LEU:HD23	1:A:184:LEU:HA	1.72	0.48
3:F:159:LYS:HZ3	3:F:169:ARG:HE	1.61	0.48
3:F:193:LEU:O	3:F:193:LEU:HD23	2.14	0.47
3:F:139:ASN:O	3:F:140:LYS:C	2.58	0.47
3:F:229:LEU:CD2	2:I:132:ILE:CD1	2.92	0.47
1:A:74:ASN:OD1	1:A:77:LYS:HB3	2.15	0.47
1:A:155:PHE:O	1:A:159:LYS:HG3	2.15	0.47
1:A:114:PRO:HA	1:A:117:LYS:HE3	1.96	0.46
3:F:179:PHE:CE1	3:F:184:LEU:HD12	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:LYS:HB2	1:A:34:GLN:OE1	2.17	0.45
1:A:74:ASN:OD1	1:A:77:LYS:HB2	2.17	0.45
3:F:50:TYR:CD1	3:F:99:VAL:CG2	3.01	0.44
3:F:5:LYS:HG2	3:F:31:VAL:HG13	1.98	0.44
3:F:33:GLU:C	3:F:35:GLY:N	2.76	0.44
3:F:33:GLU:C	3:F:35:GLY:H	2.25	0.44
3:F:37:GLU:OE2	3:F:110:ASN:O	2.36	0.43
3:F:74:ASN:HB2	3:F:77:LYS:CB	2.47	0.43
3:F:181:TYR:HB2	3:F:190:ALA:CB	2.48	0.43
2:I:135:VAL:O	2:I:136:GLN:C	2.61	0.43
3:F:114:PRO:O	3:F:118:VAL:HG23	2.19	0.43
4:F:301:OQE:C08	4:F:301:OQE:O04	2.67	0.43
2:B:135:VAL:O	2:B:135:VAL:HG12	2.18	0.42
3:F:110:ASN:O	3:F:112:THR:HG23	2.19	0.42
1:A:50:TYR:CD1	1:A:99:VAL:CG2	3.02	0.42
1:A:209:ASN:O	1:A:210:GLU:C	2.62	0.42
2:I:132:ILE:O	2:I:136:GLN:HA	2.19	0.42
3:F:36:HIS:H	3:F:36:HIS:HD1	1.67	0.42
3:F:74:ASN:HB2	3:F:77:LYS:HB3	2.00	0.42
1:A:202:ILE:C	1:A:204:GLU:H	2.27	0.42
3:F:184:LEU:O	3:F:185:ASN:C	2.63	0.42
1:A:82:LYS:HG3	1:A:85:ARG:NH1	2.34	0.41
3:F:58:ARG:HH12	4:F:301:OQE:C05	2.32	0.41
2:I:122:LEU:O	2:I:126:ILE:HD12	2.20	0.41
1:A:191:CYS:O	1:A:227:LEU:HD21	2.20	0.41
3:F:61:TRP:CD1	3:F:89:GLU:HG3	2.55	0.41
1:A:114:PRO:O	1:A:118:VAL:HG23	2.21	0.41
3:F:8:LEU:HD13	3:F:30:ALA:HB3	2.01	0.41
2:B:118:ASP:OD1	2:B:120:ILE:HD12	2.21	0.41
1:A:111:ALA:HB3	1:A:117:LYS:HG2	2.03	0.40
3:F:121:LEU:HB3	3:F:154:ALA:HB2	2.03	0.40
3:F:181:TYR:CZ	3:F:230:TRP:CD1	3.09	0.40
3:F:63:VAL:O	3:F:67:ILE:HG13	2.22	0.40
3:F:166:HIS:HE1	3:F:168:ILE:HD12	1.86	0.40
1:A:70:LYS:O	1:A:71:THR:C	2.64	0.40
1:A:128:PHE:HB3	1:A:147:SER:HB2	2.01	0.40
2:B:128:ARG:HG2	2:B:128:ARG:H	1.72	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ARG:HH22	3:F:210:GLU:OE2[4_556]	1.39	0.21
1:A:113:GLN:NE2	3:F:4:ASP:OD1[3_555]	2.07	0.13
1:A:224:ARG:NH2	3:F:210:GLU:OE2[4_556]	2.17	0.03

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	223/229 (97%)	210 (94%)	10 (4%)	3 (1%)	9	3
2	B	18/21 (86%)	16 (89%)	1 (6%)	1 (6%)	1	0
2	I	18/21 (86%)	16 (89%)	2 (11%)	0	100	100
3	F	224/231 (97%)	207 (92%)	15 (7%)	2 (1%)	14	7
All	All	483/502 (96%)	449 (93%)	28 (6%)	6 (1%)	10	4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	SER
3	F	111	ALA
3	F	185	ASN
1	A	203	ALA
1	A	206	ASP
2	B	135	VAL

5.3.2 Protein sidechains [i](#)

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	201/203 (99%)	201 (100%)	0	100	100
2	B	18/19 (95%)	18 (100%)	0	100	100
2	I	18/19 (95%)	18 (100%)	0	100	100
3	F	202/205 (98%)	202 (100%)	0	100	100
All	All	439/446 (98%)	439 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	124	ASN
3	F	69	GLN
3	F	139	ASN
3	F	221	GLN
3	F	226	ASN
2	I	124	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	$\# Z > 2$
4	OQE	F	301	-	24,24,24	1.94	3 (12%)	31,32,32	1.69	8 (25%)
4	OQE	A	301	-	24,24,24	2.50	4 (16%)	31,32,32	2.04	11 (35%)

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
4	OQE	F	301	-	-	8/17/17/17	0/2/2/2
4	OQE	A	301	-	-	7/17/17/17	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	OQE	P02-C05	8.49	1.90	1.80
4	F	301	OQE	C09-N10	5.64	1.46	1.33
4	A	301	OQE	C09-N10	5.57	1.46	1.33
4	F	301	OQE	P02-C05	5.49	1.87	1.80
4	A	301	OQE	O07-C06	3.97	1.45	1.37
4	F	301	OQE	O07-C06	3.67	1.44	1.37
4	A	301	OQE	C05-C06	2.96	1.44	1.40

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	OQE	O01-P02-O03	-4.71	99.46	112.49
4	A	301	OQE	O07-C06-C20	-4.61	113.88	123.95
4	A	301	OQE	O03-P02-C05	4.27	118.75	110.90
4	F	301	OQE	C08-C09-N10	3.77	122.98	116.63
4	A	301	OQE	O07-C06-C05	3.62	125.32	115.61
4	F	301	OQE	O04-P02-O03	-3.61	102.49	112.49
4	F	301	OQE	O07-C06-C05	3.00	123.65	115.61
4	F	301	OQE	O07-C06-C20	-2.83	117.76	123.95
4	F	301	OQE	O04-P02-C05	2.77	111.40	106.64
4	A	301	OQE	P02-C05-C06	2.59	127.15	122.15
4	A	301	OQE	C08-C09-N10	2.57	120.96	116.63
4	F	301	OQE	O01-P02-C05	2.52	110.98	106.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	301	OQE	O19-C09-N10	-2.48	118.15	123.03
4	A	301	OQE	O04-P02-O03	-2.44	105.73	112.49
4	A	301	OQE	O19-C09-N10	-2.38	118.36	123.03
4	A	301	OQE	O04-P02-C05	2.35	110.68	106.64
4	A	301	OQE	C11-C12-C13	2.24	117.89	112.83
4	A	301	OQE	O01-P02-C05	2.16	110.36	106.64
4	F	301	OQE	C22-C21-C20	-2.02	117.75	120.24

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	301	OQE	C05-C06-O07-C08
4	A	301	OQE	C08-C09-N10-C11
4	F	301	OQE	C08-C09-N10-C11
4	A	301	OQE	O19-C09-N10-C11
4	F	301	OQE	O19-C09-N10-C11
4	A	301	OQE	N10-C11-C12-C13
4	F	301	OQE	C05-C06-O07-C08
4	A	301	OQE	C20-C06-O07-C08
4	F	301	OQE	C20-C06-O07-C08
4	F	301	OQE	O07-C08-C09-O19
4	A	301	OQE	O07-C08-C09-O19
4	F	301	OQE	O07-C08-C09-N10
4	F	301	OQE	C11-C12-C13-C14
4	A	301	OQE	O07-C08-C09-N10
4	F	301	OQE	C11-C12-C13-C18

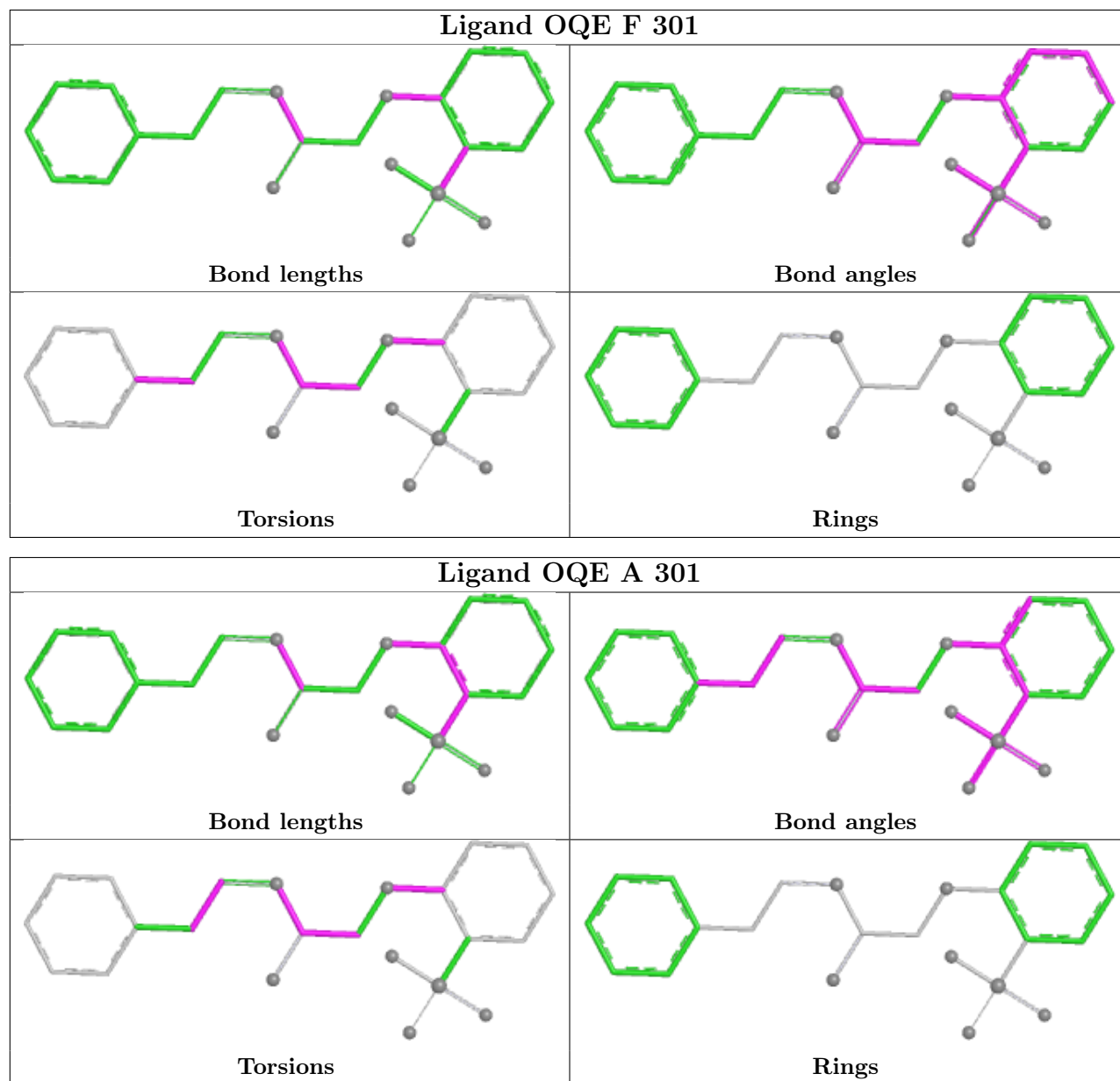
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	301	OQE	2	0
4	A	301	OQE	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	227/229 (99%)	0.48	7 (3%) 51 52	47, 73, 108, 125	0
2	B	20/21 (95%)	0.44	0 100 100	59, 71, 105, 108	0
2	I	20/21 (95%)	0.80	1 (5%) 34 34	63, 79, 108, 128	0
3	F	228/231 (98%)	0.61	14 (6%) 27 27	47, 76, 110, 130	0
All	All	495/502 (98%)	0.55	22 (4%) 39 39	47, 75, 110, 130	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	94	ASP	6.5
2	I	135	VAL	4.7
3	F	205	LEU	4.0
3	F	36	HIS	3.2
3	F	139	ASN	3.1
3	F	74	ASN	2.7
3	F	80	MET	2.6
1	A	71	THR	2.6
3	F	5	LYS	2.5
3	F	2	THR	2.5
1	A	185	ASN	2.5
1	A	136	SER	2.4
3	F	9	VAL	2.4
3	F	92	LEU	2.3
3	F	140	LYS	2.3
3	F	31	VAL	2.3
1	A	10	GLN	2.2
3	F	99	VAL	2.2
3	F	135	ALA	2.2
1	A	74	ASN	2.2
1	A	77	LYS	2.1
3	F	232	SER	2.1

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](#)

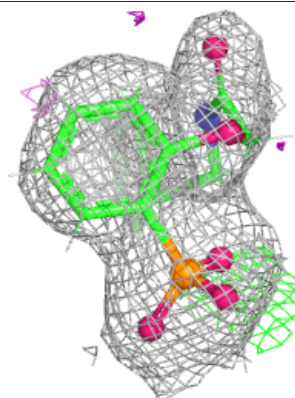
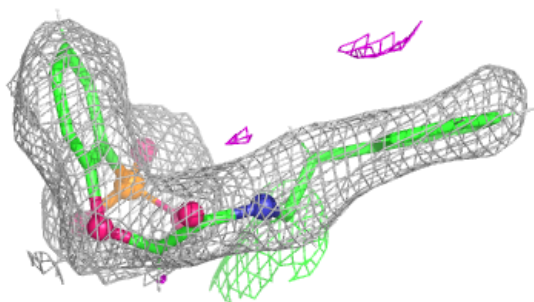
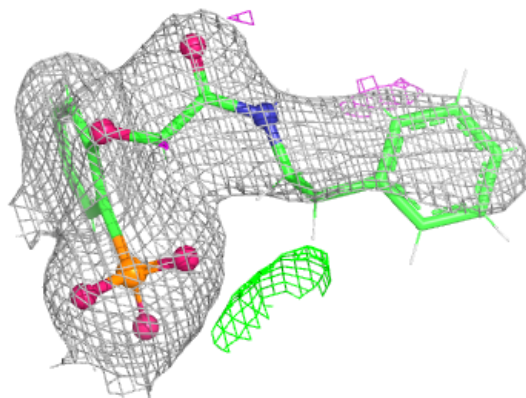
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	OQE	A	301	23/23	0.94	0.12	49,77,103,107	0
4	OQE	F	301	23/23	0.95	0.10	53,75,95,96	0

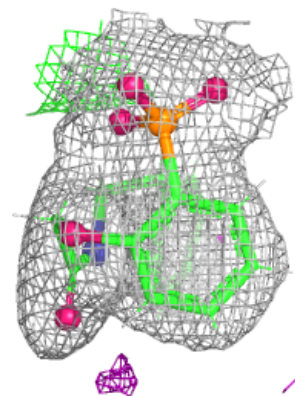
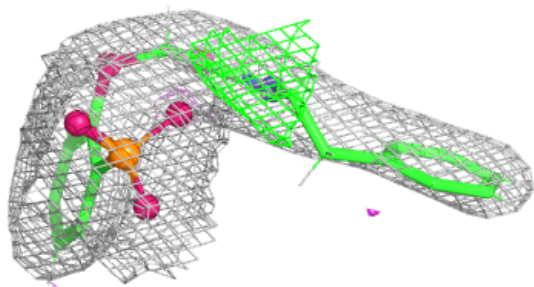
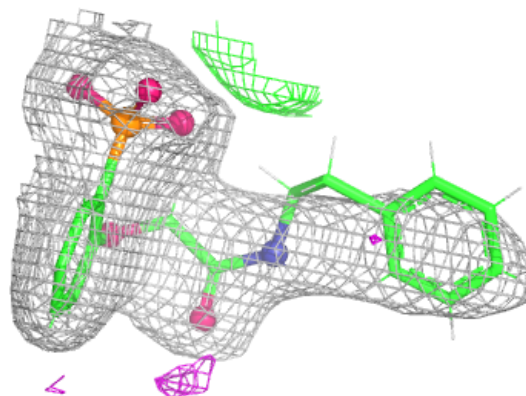
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OQE A 301:

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

**Electron density around OQE F 301:**

$2mF_o-DF_c$ (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.