



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:49 AM UTC

PDB ID : 3KS9 / pdb_00003ks9
Title : Metabotropic glutamate receptor mGluR1 complexed with LY341495 antagonist
Authors : Dobrovetsky, E.; Khutoreskaya, G.; Seitova, A.; Cossar, D.; Edwards, A.M.; Arrowsmith, C.H.; Bountra, C.; Weigelt, J.; Bochkarev, A.; Structural Genomics Consortium (SGC)
Deposited on : 2009-11-20
Resolution : 1.90 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 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

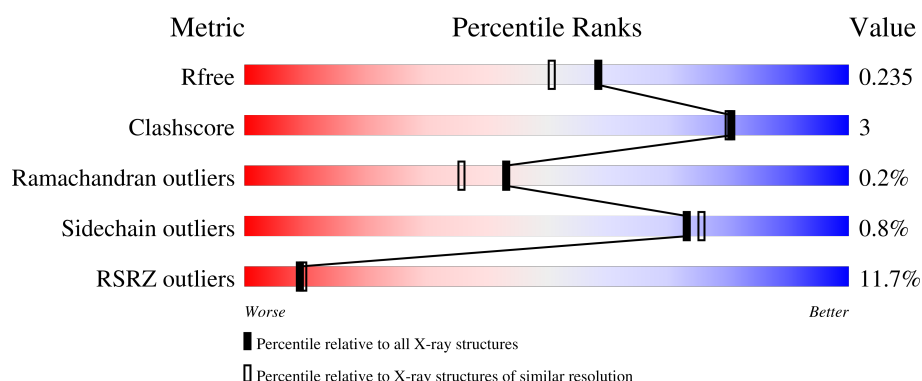
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	496	6% 83% 7% 10%
1	B	496	15% 82% 7% 11%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7727 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.

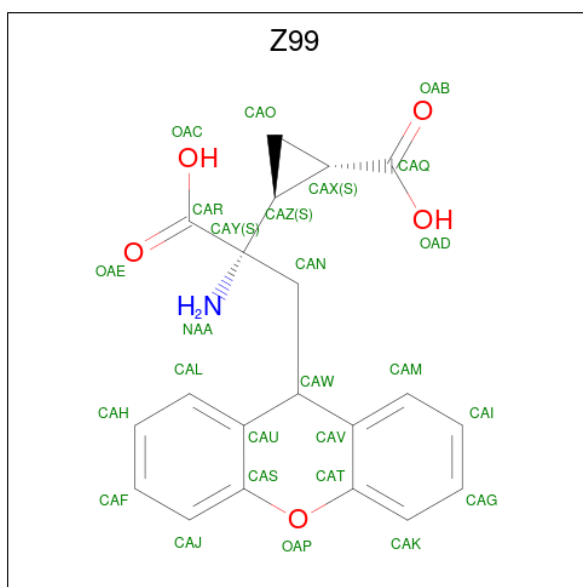
- Molecule 1 is a protein called Metabotropic glutamate receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	445	Total	C	N	O	S	0	7	0
			3565	2266	614	665	20			
1	B	442	Total	C	N	O	S	0	1	0
			3429	2182	587	641	19			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	254	SER	CYS	engineered mutation	UNP Q13255
A	519	LEU	-	expression tag	UNP Q13255
A	520	VAL	-	expression tag	UNP Q13255
A	521	PRO	-	expression tag	UNP Q13255
A	522	ARG	-	expression tag	UNP Q13255
A	523	GLY	-	expression tag	UNP Q13255
B	254	SER	CYS	engineered mutation	UNP Q13255
B	519	LEU	-	expression tag	UNP Q13255
B	520	VAL	-	expression tag	UNP Q13255
B	521	PRO	-	expression tag	UNP Q13255
B	522	ARG	-	expression tag	UNP Q13255
B	523	GLY	-	expression tag	UNP Q13255

- Molecule 2 is 2-[(1S,2S)-2-carboxycyclopropyl]-3-(9H-xanthen-9-yl)-D-alanine (CCD ID: Z99) (formula: C₂₀H₁₉NO₅).

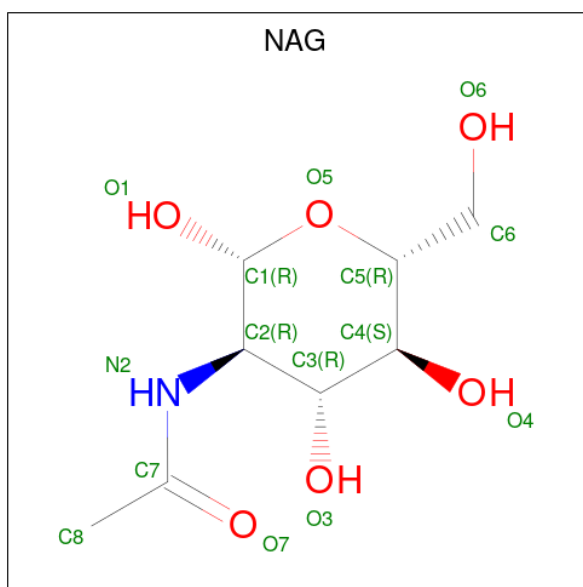


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			26	20	1	5		
2	B	1	Total	C	N	O	0	0
			26	20	1	5		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

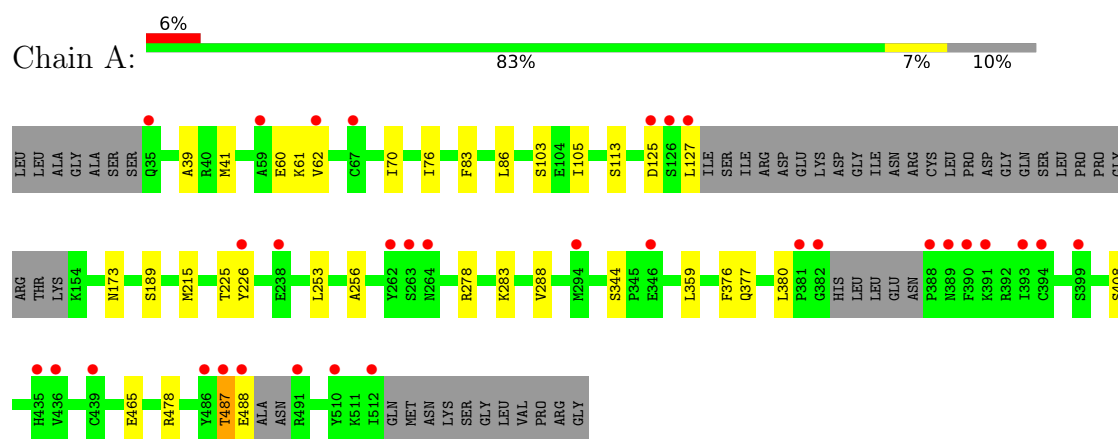
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	378	Total	O	0	0
			378	378		
5	B	273	Total	O	0	0
			273	273		

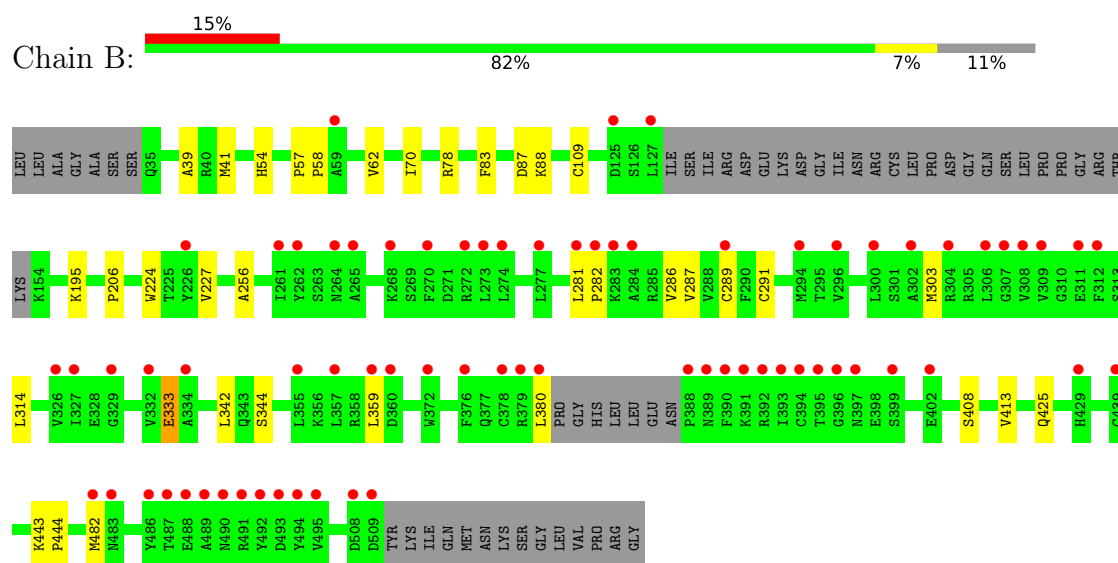
3 Residue-property plots [i](#)

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: Metabotropic glutamate receptor 1



• Molecule 1: Metabotropic glutamate receptor 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.23Å 96.55Å 97.55Å 90.00° 106.67° 90.00°	Depositor
Resolution (Å)	93.44 – 1.90 93.44 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (93.44-1.90) 98.9 (93.44-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.206 , 0.241 (Not available) , 0.235	Depositor DCC
R_{free} test set	5456 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.9	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.6	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	7727	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: MG, NAG, Z99

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	0.72	1/3645 (0.0%)	0.81	0/4930
1	B	0.66	1/3504 (0.0%)	0.78	0/4747
All	All	0.69	2/7149 (0.0%)	0.80	0/9677

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	62	VAL	CA-CB	7.13	1.57	1.54
1	A	62	VAL	CA-CB	6.04	1.57	1.54

There are no bond angle outliers.

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	3565	0	3441	17	0
1	B	3429	0	3291	18	0
2	A	26	0	17	1	0
2	B	26	0	17	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	28	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	378	0	0	2	0
5	B	273	0	0	1	0
All	All	7727	0	6792	36	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 3.

All (36) 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:173:ASN:ND2	5:A:534:HOH:O	2.42	0.53
1:A:487:THR:O	1:A:488:GLU:CB	2.58	0.51
1:B:87:ASP:OD2	5:B:712:HOH:O	2.19	0.51
1:A:215:MET:HE2	1:A:288:VAL:HG13	1.92	0.50
1:B:344:SER:HB2	1:B:408:SER:O	2.10	0.50
1:A:41:MET:HE3	1:A:83:PHE:HD2	1.77	0.49
1:B:359:LEU:HD21	1:B:380:LEU:HD21	1.95	0.48
1:B:224:TRP:CD2	1:B:286:VAL:HG21	2.49	0.47
1:A:39:ALA:HB3	1:A:105:ILE:HB	1.96	0.46
1:B:303:MET:HE2	1:B:333:GLU:HB2	1.97	0.46
1:A:189:SER:HB2	2:A:1:Z99:HAH	1.98	0.46
1:B:41:MET:HE3	1:B:83:PHE:HD2	1.81	0.45
1:B:287:VAL:HB	1:B:314:LEU:HD23	1.99	0.45
2:B:1:Z99:OAC	2:B:1:Z99:HAOA	2.16	0.44
1:B:227:VAL:HG12	1:B:286:VAL:HB	1.98	0.44
1:A:60:GLU:HG2	1:A:61:LYS:HG2	2.00	0.44
1:A:125:ASP:C	1:A:127:LEU:H	2.26	0.43
1:A:344:SER:HB2	1:A:408:SER:O	2.19	0.43
1:B:88:LYS:HE3	1:B:88:LYS:HB2	1.86	0.42
1:A:278:ARG:NH1	5:A:779:HOH:O	2.49	0.42
1:B:195:LYS:HA	1:B:195:LYS:HE2	2.02	0.42
1:B:78:ARG:HG2	1:B:413:VAL:HG21	2.01	0.42
1:A:376:PHE:O	1:A:377:GLN:C	2.63	0.42
1:B:443:LYS:HA	1:B:444:PRO:HA	1.84	0.42
1:B:289:CYS:HB3	1:B:291:CYS:SG	2.60	0.42
1:A:359:LEU:HD21	1:A:380:LEU:HD21	2.01	0.41
1:A:465:GLU:HG3	1:A:478:ARG:NH2	2.35	0.41
1:A:226[B]:TYR:CD2	1:A:283:LYS:HG2	2.55	0.41
1:B:39:ALA:HB1	1:B:83:PHE:CZ	2.56	0.41
1:B:54:HIS:O	1:B:109:CYS:HA	2.20	0.41
1:B:57:PRO:HA	1:B:58:PRO:HD2	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:PRO:HG2	1:B:342:LEU:HD21	2.03	0.41
1:B:281:LEU:HB3	1:B:282:PRO:HA	2.03	0.41
1:A:225:THR:C	1:A:253:LEU:HD12	2.47	0.40
1:A:70:ILE:HD11	1:A:76:ILE:HD11	2.03	0.40
1:A:86:LEU:CD1	1:A:103:SER:HB3	2.51	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	444/496 (90%)	432 (97%)	11 (2%)	1 (0%)	43	36
1	B	437/496 (88%)	422 (97%)	14 (3%)	1 (0%)	43	36
All	All	881/992 (89%)	854 (97%)	25 (3%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	256	ALA
1	A	256	ALA

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	378/424 (89%)	376 (100%)	2 (0%)	81	84
1	B	357/424 (84%)	353 (99%)	4 (1%)	65	67
All	All	735/848 (87%)	729 (99%)	6 (1%)	73	75

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	SER
1	A	487	THR
1	B	70	ILE
1	B	333	GLU
1	B	425	GLN
1	B	482	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	170	GLN
1	A	173	ASN
1	A	250	GLN
1	A	374	HIS
1	A	429	HIS
1	B	84	HIS
1	B	173	ASN
1	B	250	GLN
1	B	335	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

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	526	1	14,14,15	0.53	0	17,19,21	0.73	0
2	Z99	A	1	-	26,29,29	1.14	1 (3%)	34,44,44	1.42	5 (14%)
2	Z99	B	1	-	26,29,29	1.30	3 (11%)	34,44,44	1.41	4 (11%)
4	NAG	A	525	1	14,14,15	0.56	0	17,19,21	1.44	1 (5%)

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	NAG	A	526	1	-	0/6/23/26	0/1/1/1
2	Z99	A	1	-	-	1/16/38/38	0/4/4/4
2	Z99	B	1	-	-	4/16/38/38	0/4/4/4
4	NAG	A	525	1	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	Z99	CAU-CAW	2.88	1.56	1.52
2	B	1	Z99	CAN-CAY	2.61	1.57	1.54
2	A	1	Z99	CAX-CAQ	2.29	1.55	1.51
2	B	1	Z99	CAO-CAZ	2.20	1.54	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	Z99	CAO-CAZ-CAY	-4.22	115.50	122.03
4	A	525	NAG	C4-C3-C2	3.85	116.66	111.02
2	A	1	Z99	CAM-CAV-CAT	3.70	121.80	117.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	Z99	CAO-CAX-CAQ	-3.17	114.04	118.94
2	A	1	Z99	OAB-CAQ-CAX	-2.91	116.15	123.01
2	B	1	Z99	CAM-CAV-CAT	2.84	120.86	117.74
2	A	1	Z99	CAL-CAU-CAS	2.61	120.61	117.74
2	A	1	Z99	CAO-CAX-CAQ	-2.28	115.42	118.94
2	B	1	Z99	OAE-CAR-CAY	-2.23	116.78	122.53
2	A	1	Z99	CAK-CAT-CAV	-2.04	118.30	121.14

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	Z99	OAB-CAQ-CAX-CAO
2	B	1	Z99	OAD-CAQ-CAX-CAO
2	B	1	Z99	CAR-CAY-CAZ-CAO
2	A	1	Z99	OAB-CAQ-CAX-CAZ
2	B	1	Z99	OAB-CAQ-CAX-CAZ

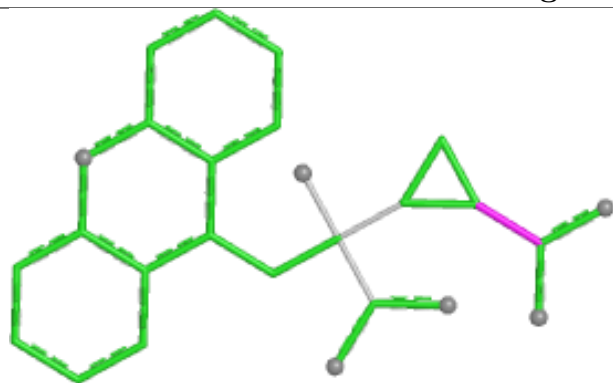
There are no ring outliers.

2 monomers are involved in 2 short contacts:

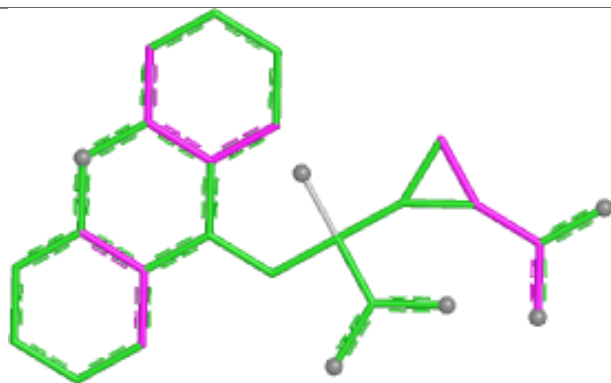
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	Z99	1	0
2	B	1	Z99	1	0

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

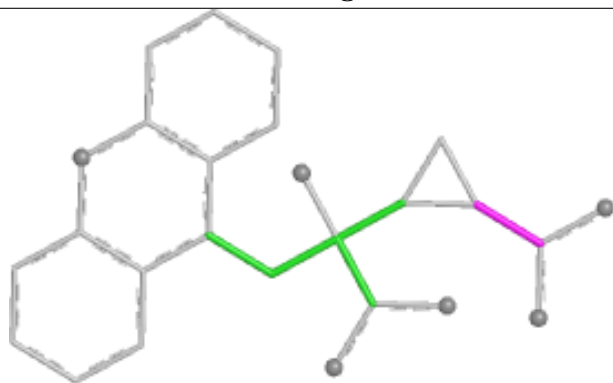
Ligand Z99 A 1



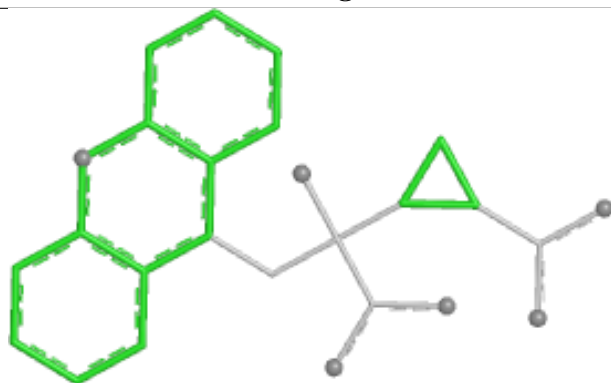
Bond lengths



Bond angles

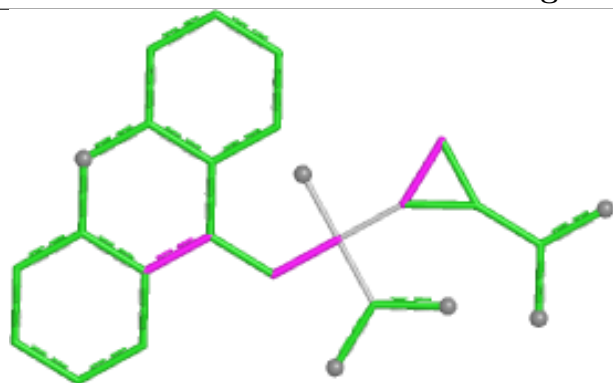


Torsions

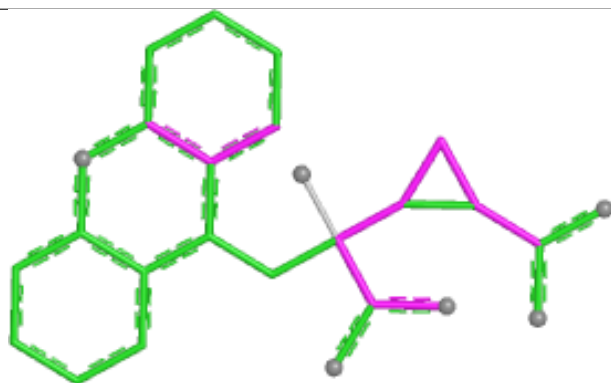


Rings

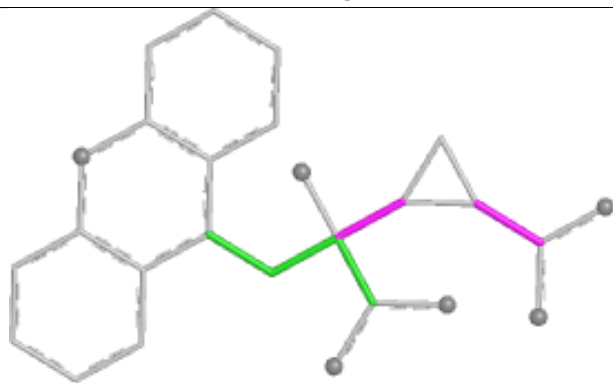
Ligand Z99 B 1



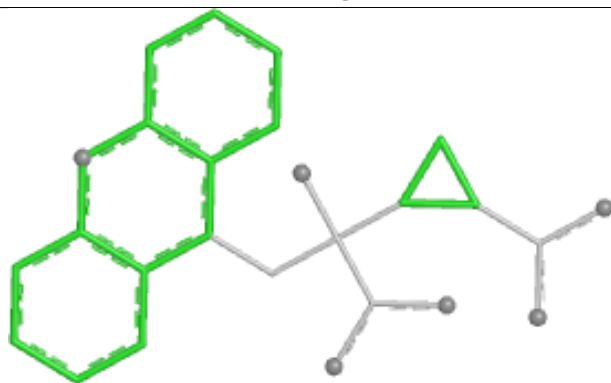
Bond lengths



Bond angles



Torsions



Rings

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	445/496 (89%)	0.39	32 (7%) 21 23	10, 30, 54, 72	7 (1%)
1	B	442/496 (89%)	0.94	72 (16%) 4 4	15, 39, 62, 76	1 (0%)
All	All	887/992 (89%)	0.67	104 (11%) 9 10	10, 34, 61, 76	8 (0%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	388	PRO	6.8
1	B	388	PRO	6.1
1	B	509	ASP	5.5
1	B	127	LEU	5.4
1	B	261	ILE	5.3
1	B	491	ARG	5.3
1	B	393	ILE	5.0
1	B	125	ASP	4.9
1	B	392	ARG	4.8
1	B	394	CYS	4.7
1	B	262	TYR	4.5
1	B	390	PHE	4.4
1	A	382	GLY	4.3
1	B	492	TYR	4.3
1	A	488	GLU	4.1
1	A	390	PHE	4.0
1	B	329	GLY	3.8
1	A	125	ASP	3.8
1	A	294[A]	MET	3.6
1	B	391	LYS	3.6
1	A	436	VAL	3.6
1	B	308	VAL	3.6
1	B	265	ALA	3.6
1	A	394	CYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	126	SER	3.5
1	B	59	ALA	3.4
1	B	508	ASP	3.4
1	A	263	SER	3.4
1	A	389	ASN	3.4
1	A	435[A]	HIS	3.4
1	B	281	LEU	3.2
1	B	380	LEU	3.2
1	A	393	ILE	3.2
1	B	264	ASN	3.1
1	B	312	PHE	3.1
1	B	376	PHE	3.1
1	B	486	TYR	3.0
1	B	283	LYS	3.0
1	B	487	THR	3.0
1	A	35	GLN	2.9
1	B	493	ASP	2.9
1	A	439	CYS	2.9
1	B	389	ASN	2.8
1	B	396	GLY	2.8
1	A	486	TYR	2.8
1	B	399	SER	2.8
1	B	372	TRP	2.8
1	B	332	VAL	2.8
1	B	327	ILE	2.8
1	B	282	PRO	2.7
1	A	226[A]	TYR	2.7
1	A	381	PRO	2.7
1	A	491	ARG	2.7
1	B	490	ASN	2.7
1	B	494	TYR	2.7
1	A	391	LYS	2.7
1	B	296	VAL	2.6
1	B	439	CYS	2.6
1	B	300	LEU	2.6
1	A	346	GLU	2.6
1	B	309	VAL	2.6
1	B	489	ALA	2.6
1	B	270	PHE	2.5
1	B	482	MET	2.5
1	B	274	LEU	2.5
1	B	359	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	127	LEU	2.5
1	B	495	VAL	2.5
1	B	378	CYS	2.4
1	A	264	ASN	2.4
1	B	429	HIS	2.4
1	B	268	LYS	2.4
1	A	67	CYS	2.4
1	B	402	GLU	2.3
1	B	397	ASN	2.3
1	B	306	LEU	2.3
1	B	307	GLY	2.3
1	A	62	VAL	2.3
1	B	488	GLU	2.3
1	B	302	ALA	2.3
1	B	273	LEU	2.3
1	B	277	LEU	2.3
1	B	355	LEU	2.3
1	B	226	TYR	2.2
1	B	379	ARG	2.2
1	A	399	SER	2.2
1	A	59	ALA	2.2
1	B	326	VAL	2.2
1	B	395	THR	2.2
1	A	262	TYR	2.1
1	A	487	THR	2.1
1	B	284	ALA	2.1
1	B	360	ASP	2.1
1	A	238	GLU	2.1
1	B	289	CYS	2.1
1	B	334	ALA	2.1
1	B	304	ARG	2.1
1	B	311	GLU	2.1
1	B	483	ASN	2.1
1	B	357	LEU	2.1
1	B	272	ARG	2.1
1	A	512	ILE	2.0
1	A	510	TYR	2.0
1	B	294	MET	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

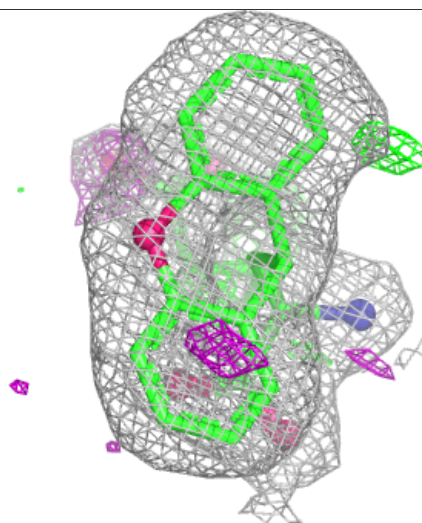
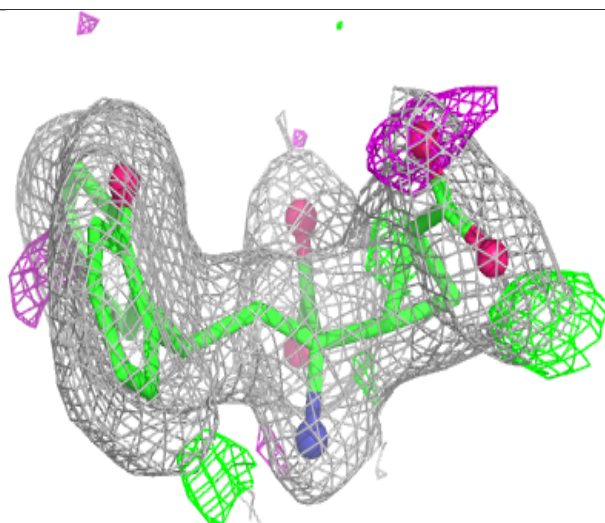
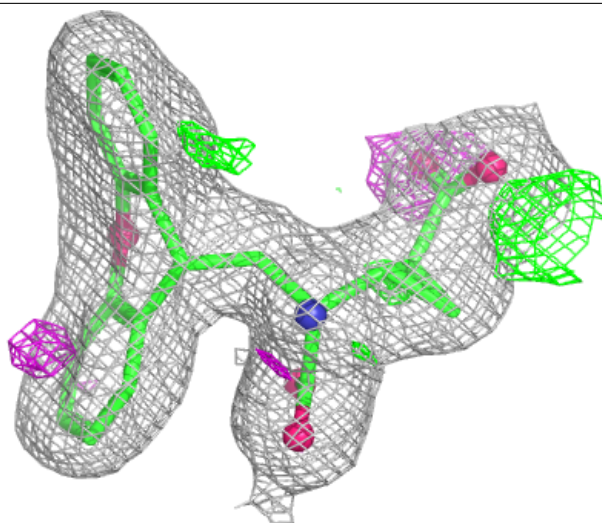
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NAG	A	526	14/15	0.62	0.14	72,74,76,76	0
4	NAG	A	525	14/15	0.77	0.14	57,62,63,64	0
2	Z99	A	1	26/26	0.92	0.09	24,28,33,36	0
2	Z99	B	1	26/26	0.93	0.09	24,31,42,45	0
3	MG	A	524	1/1	0.98	0.11	21,21,21,21	0
3	MG	B	524	1/1	0.98	0.07	26,26,26,26	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.

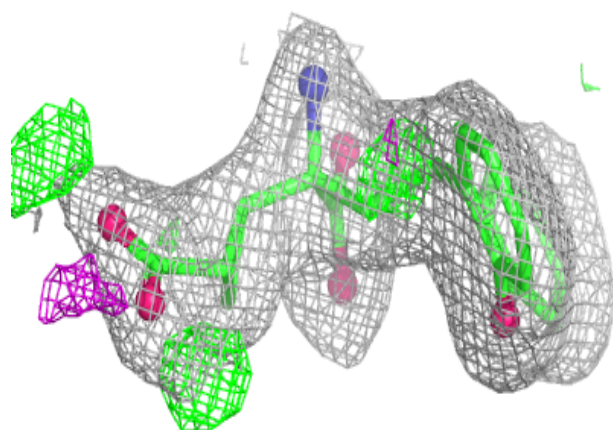
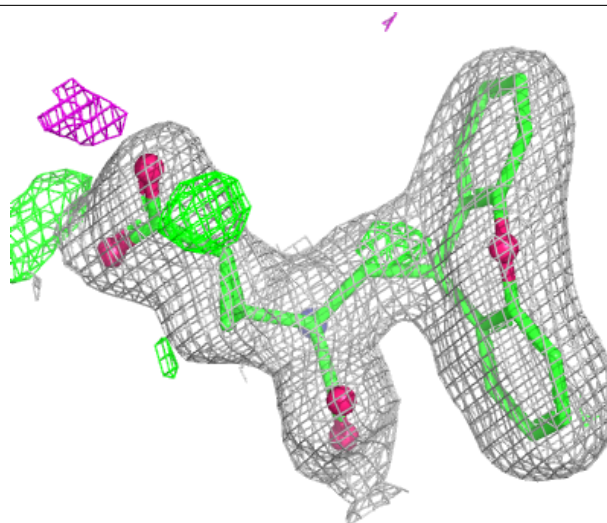
Electron density around Z99 A 1:

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



Electron density around Z99 B 1:

$2mF_o-DF_c$ (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.