Full wwPDB X-ray Structure Validation Report (i)

Dec 17, 2023 – 03:57 PM EST

PDB ID	:	4U5B
Title	:	Crystal structure of GluA2 A622T, con-ikot-ikot snail toxin, partial agonist KA and postitive modulator (R,R)-2b complex
Authors	:	Chen, L.; Gouaux, E.
Deposited on	:	2014-07-25
Resolution	:	3.50 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 (i) 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 (i)) were used in the production of this report:

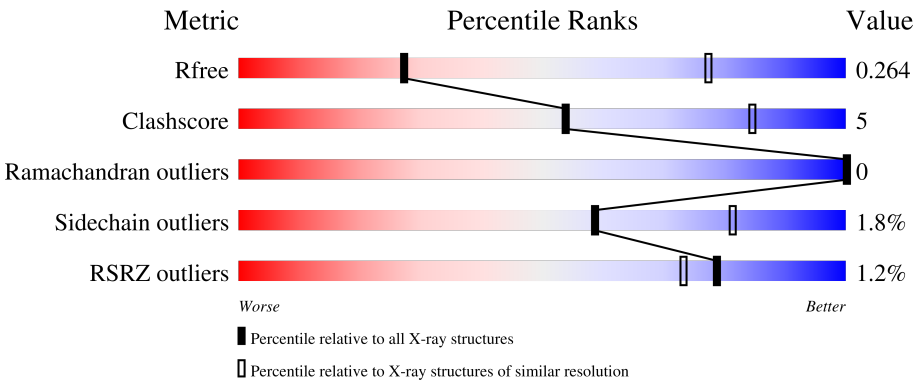
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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 <=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	814	% 78% 12% • 9%
1	B	814	% 78% 13% 9%
1	C	814	% 80% 11% • 8%
1	D	814	% 78% 12% 9%
2	E	90	79% 16% 6%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	90	79% 16% 6%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 23751 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 Glutamate receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	738	5558	3578	905	1049	26	0	0	0
1	B	743	5646	3630	929	1061	26	0	0	0
1	C	746	5507	3534	903	1044	26	0	0	0
1	D	738	5578	3594	906	1052	26	0	0	0

There are 124 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	184	GLY	LYS	engineered mutation	UNP P19491
A	237	GLU	ASN	engineered mutation	UNP P19491
A	?	-	LEU	deletion	UNP P19491
A	?	-	PRO	deletion	UNP P19491
A	385	ASP	ASN	engineered mutation	UNP P19491
A	392	GLN	ASN	engineered mutation	UNP P19491
A	461	ASP	ASN	engineered mutation	UNP P19491
A	528	ALA	CYS	engineered mutation	UNP P19491
A	535	LEU	GLY	engineered mutation	UNP P19491
A	?	-	ARG	deletion	UNP P19491
A	?	-	GLU	deletion	UNP P19491
A	?	-	THR	deletion	UNP P19491
A	?	-	GLN	deletion	UNP P19491
A	?	-	SER	deletion	UNP P19491
A	565	GLU	SER	engineered mutation	UNP P19491
A	577	PHE	LEU	engineered mutation	UNP P19491
A	580	ALA	SER	engineered mutation	UNP P19491
A	582	LYS	GLY	engineered mutation	UNP P19491
A	583	LEU	ALA	engineered mutation	UNP P19491
A	585	PHE	MET	engineered mutation	UNP P19491
A	589	ALA	CYS	engineered mutation	UNP P19491

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	598	ALA	GLY	engineered mutation	UNP P19491
A	602	ALA	GLY	engineered mutation	UNP P19491
A	622	THR	ALA	engineered mutation	UNP P19491
A	815	ALA	CYS	engineered mutation	UNP P19491
A	818	ARG	SER	engineered mutation	UNP P19491
A	819	MET	ARG	engineered mutation	UNP P19491
A	820	LYS	ALA	engineered mutation	UNP P19491
A	821	LEU	GLU	engineered mutation	UNP P19491
A	822	VAL	ALA	engineered mutation	UNP P19491
A	823	PRO	LYS	engineered mutation	UNP P19491
B	184	GLY	LYS	engineered mutation	UNP P19491
B	237	GLU	ASN	engineered mutation	UNP P19491
B	?	-	LEU	deletion	UNP P19491
B	?	-	PRO	deletion	UNP P19491
B	385	ASP	ASN	engineered mutation	UNP P19491
B	392	GLN	ASN	engineered mutation	UNP P19491
B	461	ASP	ASN	engineered mutation	UNP P19491
B	528	ALA	CYS	engineered mutation	UNP P19491
B	535	LEU	GLY	engineered mutation	UNP P19491
B	?	-	ARG	deletion	UNP P19491
B	?	-	GLU	deletion	UNP P19491
B	?	-	THR	deletion	UNP P19491
B	?	-	GLN	deletion	UNP P19491
B	?	-	SER	deletion	UNP P19491
B	565	GLU	SER	engineered mutation	UNP P19491
B	577	PHE	LEU	engineered mutation	UNP P19491
B	580	ALA	SER	engineered mutation	UNP P19491
B	582	LYS	GLY	engineered mutation	UNP P19491
B	583	LEU	ALA	engineered mutation	UNP P19491
B	585	PHE	MET	engineered mutation	UNP P19491
B	589	ALA	CYS	engineered mutation	UNP P19491
B	598	ALA	GLY	engineered mutation	UNP P19491
B	602	ALA	GLY	engineered mutation	UNP P19491
B	622	THR	ALA	engineered mutation	UNP P19491
B	815	ALA	CYS	engineered mutation	UNP P19491
B	818	ARG	SER	engineered mutation	UNP P19491
B	819	MET	ARG	engineered mutation	UNP P19491
B	820	LYS	ALA	engineered mutation	UNP P19491
B	821	LEU	GLU	engineered mutation	UNP P19491
B	822	VAL	ALA	engineered mutation	UNP P19491
B	823	PRO	LYS	engineered mutation	UNP P19491
C	184	GLY	LYS	engineered mutation	UNP P19491

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	237	GLU	ASN	engineered mutation	UNP P19491
C	?	-	LEU	deletion	UNP P19491
C	?	-	PRO	deletion	UNP P19491
C	385	ASP	ASN	engineered mutation	UNP P19491
C	392	GLN	ASN	engineered mutation	UNP P19491
C	461	ASP	ASN	engineered mutation	UNP P19491
C	528	ALA	CYS	engineered mutation	UNP P19491
C	535	LEU	GLY	engineered mutation	UNP P19491
C	?	-	ARG	deletion	UNP P19491
C	?	-	GLU	deletion	UNP P19491
C	?	-	THR	deletion	UNP P19491
C	?	-	GLN	deletion	UNP P19491
C	?	-	SER	deletion	UNP P19491
C	565	GLU	SER	engineered mutation	UNP P19491
C	577	PHE	LEU	engineered mutation	UNP P19491
C	580	ALA	SER	engineered mutation	UNP P19491
C	582	LYS	GLY	engineered mutation	UNP P19491
C	583	LEU	ALA	engineered mutation	UNP P19491
C	585	PHE	MET	engineered mutation	UNP P19491
C	589	ALA	CYS	engineered mutation	UNP P19491
C	598	ALA	GLY	engineered mutation	UNP P19491
C	602	ALA	GLY	engineered mutation	UNP P19491
C	622	THR	ALA	engineered mutation	UNP P19491
C	815	ALA	CYS	engineered mutation	UNP P19491
C	818	ARG	SER	engineered mutation	UNP P19491
C	819	MET	ARG	engineered mutation	UNP P19491
C	820	LYS	ALA	engineered mutation	UNP P19491
C	821	LEU	GLU	engineered mutation	UNP P19491
C	822	VAL	ALA	engineered mutation	UNP P19491
C	823	PRO	LYS	engineered mutation	UNP P19491
D	184	GLY	LYS	engineered mutation	UNP P19491
D	237	GLU	ASN	engineered mutation	UNP P19491
D	?	-	LEU	deletion	UNP P19491
D	?	-	PRO	deletion	UNP P19491
D	385	ASP	ASN	engineered mutation	UNP P19491
D	392	GLN	ASN	engineered mutation	UNP P19491
D	461	ASP	ASN	engineered mutation	UNP P19491
D	528	ALA	CYS	engineered mutation	UNP P19491
D	535	LEU	GLY	engineered mutation	UNP P19491
D	?	-	ARG	deletion	UNP P19491
D	?	-	GLU	deletion	UNP P19491
D	?	-	THR	deletion	UNP P19491

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	GLN	deletion	UNP P19491
D	?	-	SER	deletion	UNP P19491
D	565	GLU	SER	engineered mutation	UNP P19491
D	577	PHE	LEU	engineered mutation	UNP P19491
D	580	ALA	SER	engineered mutation	UNP P19491
D	582	LYS	GLY	engineered mutation	UNP P19491
D	583	LEU	ALA	engineered mutation	UNP P19491
D	585	PHE	MET	engineered mutation	UNP P19491
D	589	ALA	CYS	engineered mutation	UNP P19491
D	598	ALA	GLY	engineered mutation	UNP P19491
D	602	ALA	GLY	engineered mutation	UNP P19491
D	622	THR	ALA	engineered mutation	UNP P19491
D	815	ALA	CYS	engineered mutation	UNP P19491
D	818	ARG	SER	engineered mutation	UNP P19491
D	819	MET	ARG	engineered mutation	UNP P19491
D	820	LYS	ALA	engineered mutation	UNP P19491
D	821	LEU	GLU	engineered mutation	UNP P19491
D	822	VAL	ALA	engineered mutation	UNP P19491
D	823	PRO	LYS	engineered mutation	UNP P19491

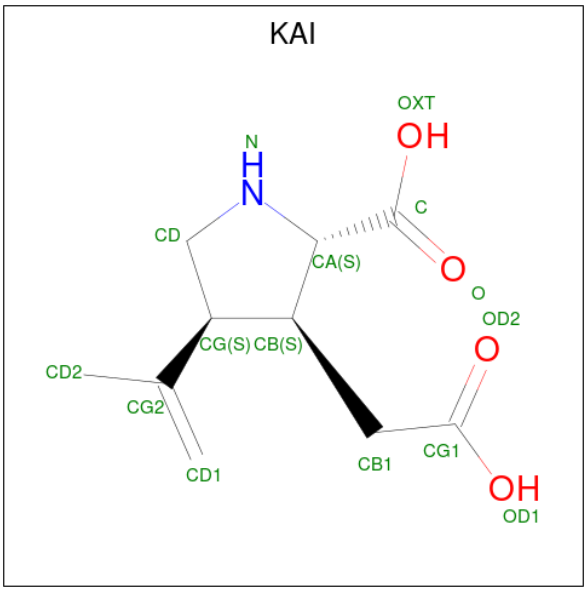
- Molecule 2 is a protein called Con-ikot-ikot.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	E	85	641	387	113	125	16	0	0	0
2	F	85	641	387	113	125	16	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

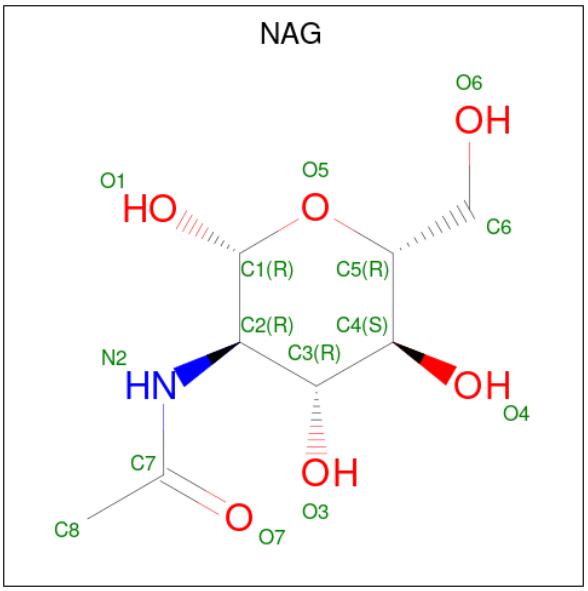
Chain	Residue	Modelled	Actual	Comment	Reference
E	-3	GLY	-	expression tag	UNP P0CB20
E	-2	PRO	-	expression tag	UNP P0CB20
E	-1	GLY	-	expression tag	UNP P0CB20
E	0	SER	-	expression tag	UNP P0CB20
F	-3	GLY	-	expression tag	UNP P0CB20
F	-2	PRO	-	expression tag	UNP P0CB20
F	-1	GLY	-	expression tag	UNP P0CB20
F	0	SER	-	expression tag	UNP P0CB20

- Molecule 3 is 3-(CARBOXYMETHYL)-4-ISOPROPENYLPROLINE (three-letter code: KAI) (formula: $C_{10}H_{15}NO_4$).



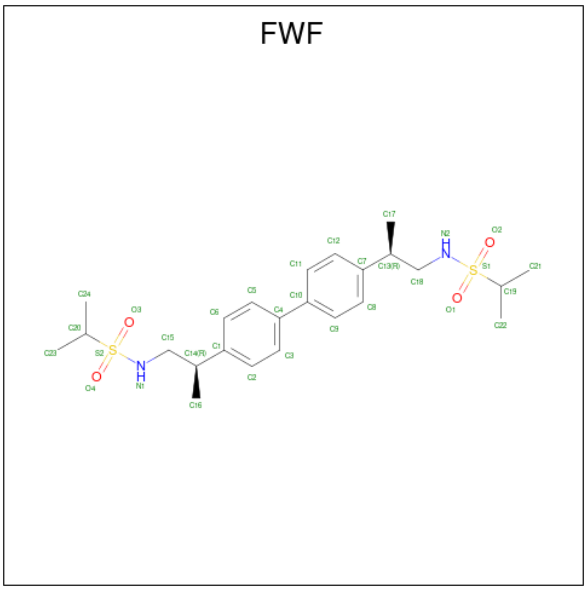
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	15	10	1	4	0	0
			Total	C	N	O		
3	B	1	15	10	1	4	0	0
			Total	C	N	O		
3	C	1	15	10	1	4	0	0
			Total	C	N	O		
3	D	1	15	10	1	4	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is N,N'-[biphenyl-4,4'-diyldi(2R)propane-2,1-diyl]dipropane-2-sulfonamide (three-letter code: FWF) (formula: $C_{24}H_{36}N_2O_4S_2$).



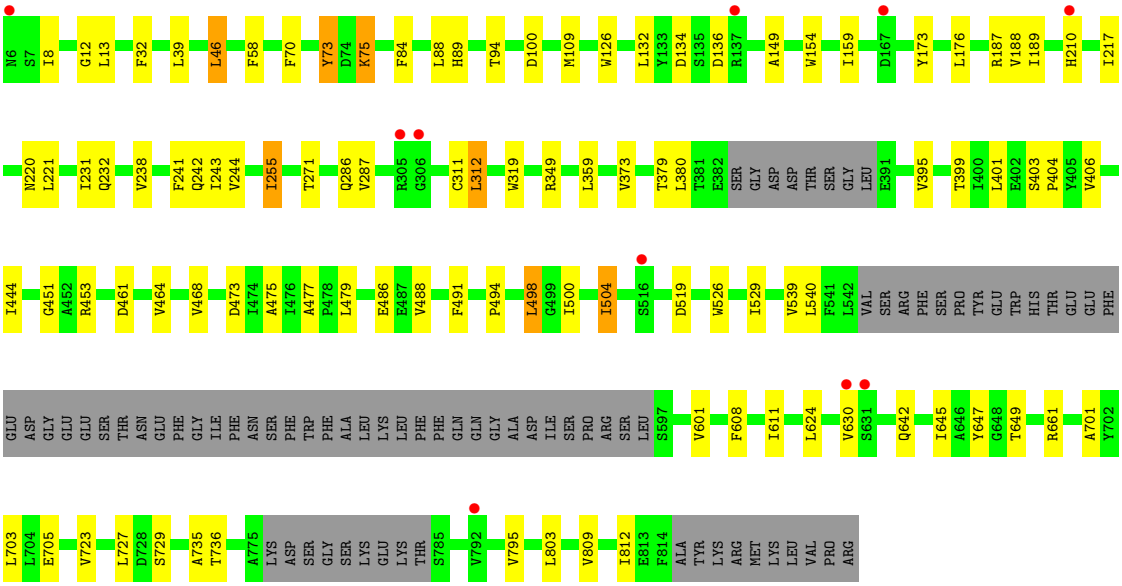
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
5	B	1	32	24	2	4	2	0	0
5	D	1	32	24	2	4	2	0	0

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.

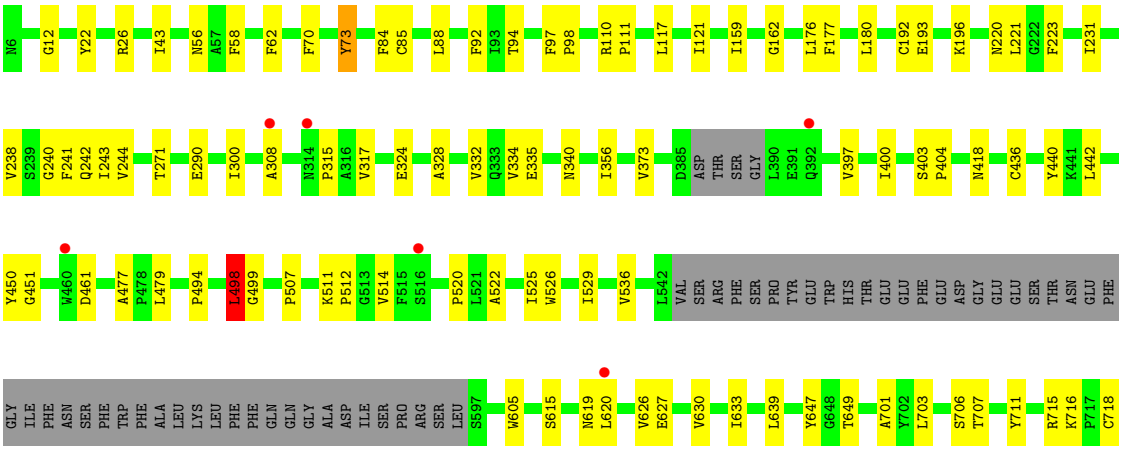
• Molecule 1: Glutamate receptor 2

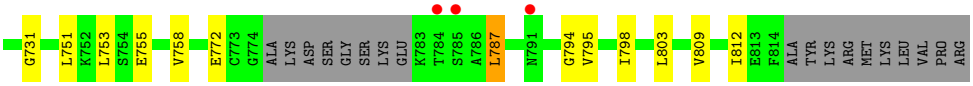
Chain A:



• Molecule 1: Glutamate receptor 2

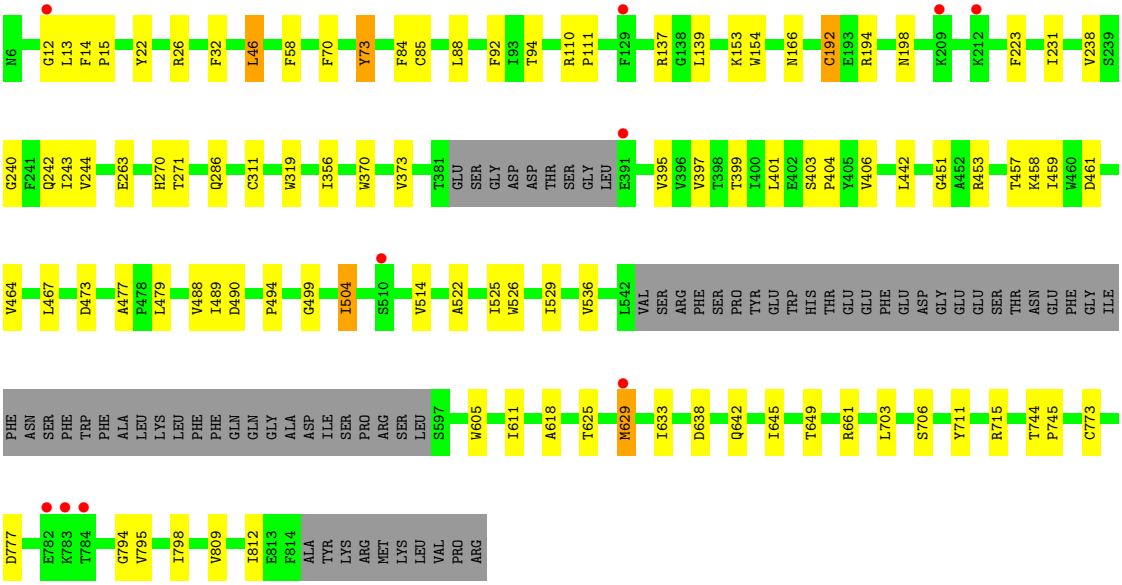
Chain B:





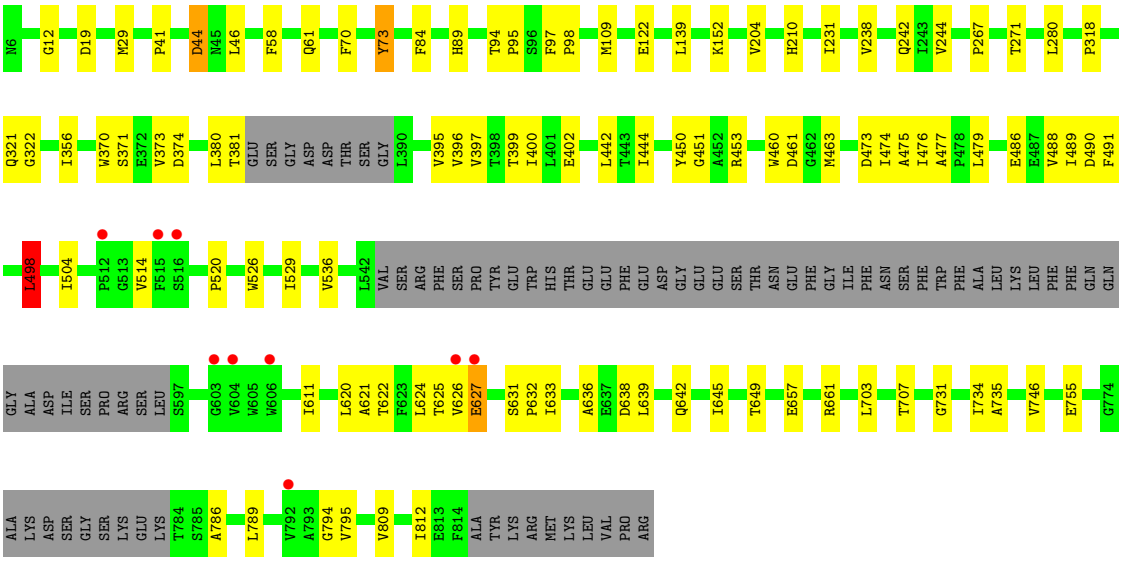
• Molecule 1: Glutamate receptor 2

Chain C:



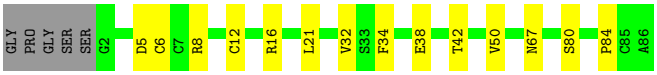
• Molecule 1: Glutamate receptor 2

Chain D:



• Molecule 2: Con-ikot-ikot

Chain E:



- Molecule 2: Con-ikot-ikot

Chain F:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	161.52Å 368.66Å 109.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 90.44 – 3.50	Depositor EDS
% Data completeness (in resolution range)	79.2 (20.00-3.50) 79.2 (90.44-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.69 (at 3.49Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.220 , 0.261 0.228 , 0.264	Depositor DCC
R_{free} test set	3305 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	103.5	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.23 , 93.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.42$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	23751	wwPDB-VP
Average B, all atoms (Å^2)	122.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: NAG, FWF, KAI

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	0.23	0/5674	0.43	0/7718
1	B	0.24	0/5762	0.43	1/7824 (0.0%)
1	C	0.23	0/5620	0.42	0/7656
1	D	0.24	0/5694	0.43	1/7747 (0.0%)
2	E	0.26	0/651	0.44	0/873
2	F	0.24	0/651	0.42	0/873
All	All	0.24	0/24052	0.42	2/32691 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	498	LEU	CA-CB-CG	5.17	127.18	115.30
1	B	498	LEU	CA-CB-CG	5.06	126.93	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5558	0	5320	60	0
1	B	5646	0	5447	66	0
1	C	5507	0	5178	55	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	5578	0	5353	61	0
2	E	641	0	593	13	0
2	F	641	0	593	10	0
3	A	15	0	13	1	0
3	B	15	0	13	0	0
3	C	15	0	13	1	0
3	D	15	0	13	2	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
4	C	14	0	13	0	0
4	D	14	0	13	0	0
5	B	32	0	36	3	0
5	D	32	0	36	2	0
All	All	23751	0	22660	247	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 (247) 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:C:464:VAL:HG13	1:C:489:ILE:HD11	1.66	0.77
1:C:649:THR:HG22	1:C:703:LEU:HB2	1.73	0.71
1:D:657:GLU:OE1	1:D:661:ARG:NH1	2.24	0.71
1:C:356:ILE:HD11	1:C:370:TRP:HB2	1.74	0.70
1:D:488:VAL:HG23	1:D:489:ILE:HG12	1.77	0.67
1:A:661:ARG:NH2	1:D:755:GLU:OE1	2.28	0.67
1:D:122:GLU:OE1	1:D:152:LYS:NZ	2.28	0.67
1:B:520:PRO:HG2	1:B:620:LEU:HD23	1.77	0.66
1:D:44:ASP:OD1	1:D:44:ASP:N	2.27	0.66
1:C:453:ARG:NH2	2:F:38:GLU:OE1	2.28	0.66
1:D:356:ILE:HD11	1:D:370:TRP:HB2	1.79	0.65
1:A:379:THR:HG22	1:A:380:LEU:H	1.62	0.64
1:B:162:GLY:HA2	1:B:196:LYS:HE2	1.78	0.64
1:B:649:THR:HG22	1:B:703:LEU:HB2	1.79	0.64
2:E:5:ASP:OD1	2:E:6:CYS:N	2.28	0.61
1:D:476:ILE:HG12	1:D:734:ILE:HD12	1.83	0.61
1:A:649:THR:HG22	1:A:703:LEU:HB2	1.81	0.61
1:B:755:GLU:OE1	1:C:661:ARG:NH2	2.34	0.60
1:A:88:LEU:HD11	1:A:311:CYS:HB2	1.82	0.60
1:D:633:ILE:HD13	1:D:639:LEU:HG	1.83	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:642:GLN:HE22	1:A:645:ILE:HB	1.67	0.59
2:F:5:ASP:OD1	2:F:6:CYS:N	2.30	0.59
1:A:232:GLN:HA	1:A:359:LEU:HD21	1.85	0.59
1:C:611:ILE:HG21	1:D:795:VAL:HG21	1.84	0.58
1:D:633:ILE:HG23	1:D:638:ASP:HB2	1.85	0.58
1:D:231:ILE:HD12	1:D:238:VAL:HG21	1.86	0.57
1:B:711:TYR:OH	1:B:715:ARG:NH1	2.38	0.56
2:E:21:LEU:HD21	2:E:32:VAL:HA	1.88	0.56
1:C:401:LEU:HD23	1:C:406:VAL:HG12	1.87	0.56
1:A:539:VAL:HG21	1:B:803:LEU:HB3	1.89	0.55
1:B:290:GLU:HG3	1:B:334:VAL:HG11	1.88	0.55
1:C:467:LEU:HD23	1:C:489:ILE:HG21	1.89	0.54
1:C:488:VAL:HG23	1:C:489:ILE:HG13	1.90	0.53
1:D:396:VAL:HG23	1:D:473:ASP:H	1.72	0.53
1:A:453:ARG:NH2	2:E:38:GLU:OE2	2.40	0.52
1:A:401:LEU:HD23	1:A:406:VAL:HG12	1.91	0.52
1:C:489:ILE:HG22	1:C:490:ASP:H	1.74	0.52
1:D:626:VAL:O	1:D:627:GLU:HG3	2.09	0.52
1:D:649:THR:HG22	1:D:703:LEU:HB2	1.92	0.52
1:C:522:ALA:HB3	1:C:525:ILE:HG13	1.90	0.52
1:C:242:GLN:HE21	1:C:244:VAL:H	1.58	0.51
1:D:109:MET:HB3	1:D:280:LEU:HD22	1.92	0.51
1:D:453:ARG:NH2	1:D:460:TRP:HE1	2.08	0.51
1:C:522:ALA:H	1:C:525:ILE:HD12	1.76	0.51
1:D:625:THR:HG21	1:D:786:ALA:CB	2.41	0.51
1:C:139:LEU:HD12	1:D:139:LEU:HD12	1.93	0.51
1:A:187:ARG:NH2	2:E:67:ASN:OD1	2.43	0.51
1:C:22:TYR:CE2	1:C:26:ARG:HD2	2.45	0.51
1:A:729:SER:HB2	5:D:903:FWF:H19	1.93	0.51
1:D:489:ILE:HG22	1:D:490:ASP:H	1.76	0.50
1:A:13:LEU:HD23	1:A:46:LEU:HD21	1.92	0.50
1:B:451:GLY:HA2	1:B:461:ASP:O	2.11	0.50
1:D:451:GLY:HA2	1:D:461:ASP:O	2.09	0.50
1:B:477:ALA:O	1:B:479:LEU:N	2.44	0.50
1:A:242:GLN:HE21	1:A:244:VAL:H	1.59	0.50
1:D:486:GLU:HG2	1:D:491:PHE:HD2	1.77	0.50
2:E:80:SER:HB2	2:F:3:PRO:HG2	1.94	0.50
1:A:608:PHE:O	1:A:611:ILE:HG13	2.10	0.49
1:A:109:MET:HE1	1:A:287:VAL:HG21	1.94	0.49
1:D:520:PRO:HG2	1:D:620:LEU:HD23	1.93	0.49
1:D:450:TYR:O	1:D:463:MET:N	2.46	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:PHE:CD1	1:C:240:GLY:HA3	2.48	0.49
1:B:514:VAL:HG13	1:B:794:GLY:HA3	1.94	0.49
1:C:73:TYR:CE2	1:C:94:THR:HG21	2.48	0.48
1:C:457:THR:O	1:C:458:LYS:HG2	2.13	0.48
1:A:611:ILE:HG12	1:B:795:VAL:HG11	1.95	0.48
1:A:312:LEU:O	1:B:56:ASN:ND2	2.41	0.48
1:B:522:ALA:HB3	1:B:525:ILE:HG13	1.94	0.48
1:A:132:LEU:HD13	1:A:159:ILE:HB	1.93	0.48
1:B:221:LEU:HB3	1:B:243:ILE:HB	1.95	0.48
1:C:88:LEU:HD11	1:C:311:CYS:HB2	1.96	0.48
1:B:315:PRO:HB2	1:B:317:VAL:O	2.14	0.48
1:C:194:ARG:O	1:C:198:ASN:ND2	2.47	0.48
1:B:400:ILE:HG21	1:B:450:TYR:CZ	2.49	0.47
1:A:488:VAL:HG12	2:E:34:PHE:HD1	1.78	0.47
1:B:809:VAL:HA	1:B:812:ILE:HG12	1.95	0.47
1:C:58:PHE:CE2	1:C:84:PHE:HB3	2.49	0.47
2:E:34:PHE:O	2:E:38:GLU:HG2	2.13	0.47
1:B:418:ASN:ND2	1:B:440:TYR:O	2.44	0.47
1:A:126:TRP:CG	1:A:187:ARG:HD3	2.49	0.47
1:B:159:ILE:HG21	1:B:176:LEU:HD13	1.96	0.47
1:A:231:ILE:HD12	1:A:238:VAL:HG21	1.95	0.47
1:B:315:PRO:C	1:B:317:VAL:H	2.18	0.47
1:D:204:VAL:HG12	1:D:210:HIS:HB3	1.96	0.47
1:D:397:VAL:HG13	1:D:474:ILE:HG23	1.95	0.47
1:A:255:ILE:HD13	1:A:255:ILE:HA	1.81	0.47
1:A:624:LEU:HB2	1:D:622:THR:HG21	1.97	0.47
1:B:494:PRO:O	5:B:903:FWF:N1	2.39	0.47
1:D:636:ALA:HA	1:D:639:LEU:HD12	1.96	0.47
1:A:73:TYR:CE2	1:A:94:THR:HG21	2.50	0.47
1:A:500:ILE:HB	1:A:727:LEU:HB2	1.95	0.47
1:C:451:GLY:HA2	1:C:461:ASP:O	2.14	0.47
1:A:8:ILE:HB	1:A:39:LEU:HD23	1.97	0.46
1:D:809:VAL:HA	1:D:812:ILE:HG12	1.97	0.46
2:F:24:TYR:O	2:F:28:GLU:HG3	2.15	0.46
1:B:177:PHE:HA	1:B:180:LEU:HB2	1.97	0.46
1:B:498:LEU:HD13	1:B:731:GLY:HA2	1.97	0.46
1:C:397:VAL:HG13	1:C:442:LEU:HA	1.97	0.46
1:A:803:LEU:HD21	1:D:536:VAL:HG22	1.97	0.46
3:C:901:KAI:HD12	3:C:901:KAI:HD2	1.69	0.46
1:A:221:LEU:HB3	1:A:243:ILE:HB	1.97	0.46
1:C:642:GLN:HE22	1:C:645:ILE:HB	1.80	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:SER:HA	1:C:404:PRO:HA	1.75	0.46
1:D:477:ALA:O	1:D:479:LEU:N	2.46	0.46
1:C:711:TYR:OH	1:C:715:ARG:NH1	2.49	0.46
1:A:58:PHE:CE2	1:A:84:PHE:HB3	2.51	0.46
1:B:397:VAL:HB	1:B:442:LEU:HD23	1.97	0.45
1:D:642:GLN:HE22	1:D:645:ILE:HB	1.82	0.45
1:B:22:TYR:CE2	1:B:26:ARG:HD2	2.52	0.45
1:B:536:VAL:HG21	1:B:605:TRP:CE3	2.50	0.45
1:B:753:LEU:HD22	1:B:758:VAL:HG11	1.97	0.45
1:D:94:THR:HA	1:D:95:PRO:HD3	1.78	0.45
1:C:231:ILE:HD12	1:C:238:VAL:HG21	1.98	0.45
1:C:514:VAL:HG13	1:C:794:GLY:HA3	1.99	0.45
1:B:62:PHE:CZ	1:B:308:ALA:HB1	2.52	0.45
1:B:633:ILE:HD13	1:B:639:LEU:HD13	1.98	0.45
1:D:475:ALA:HB3	1:D:735:ALA:HB3	1.97	0.45
1:D:498:LEU:HD13	1:D:731:GLY:HA2	1.98	0.45
1:B:231:ILE:HD12	1:B:238:VAL:HG21	1.99	0.45
1:B:300:ILE:HG21	1:B:324:GLU:HG2	1.98	0.45
1:A:486:GLU:HG2	1:A:491:PHE:HD2	1.82	0.45
1:D:29:MET:SD	1:D:41:PRO:HG3	2.56	0.45
1:D:242:GLN:HE21	1:D:244:VAL:H	1.65	0.45
2:F:18:ASN:HB2	2:F:73:HIS:CE1	2.51	0.45
1:A:100:ASP:OD1	1:A:100:ASP:N	2.46	0.45
1:B:12:GLY:HA2	1:B:70:PHE:O	2.17	0.45
1:B:223:PHE:CD1	1:B:240:GLY:HA3	2.52	0.45
1:C:13:LEU:HD23	1:C:46:LEU:HD21	1.99	0.45
1:A:395:VAL:HG13	1:A:473:ASP:HB2	1.99	0.45
1:A:451:GLY:HA2	1:A:461:ASP:O	2.17	0.45
1:B:193:GLU:H	1:B:193:GLU:HG2	1.63	0.45
1:D:73:TYR:CE2	1:D:94:THR:HG21	2.52	0.45
1:D:734:ILE:HG21	1:D:746:VAL:HG11	1.99	0.45
1:A:12:GLY:HA2	1:A:70:PHE:O	2.17	0.44
1:B:98:PRO:HD3	1:B:110:ARG:HD2	1.99	0.44
1:A:32:PHE:CZ	1:A:286:GLN:HB2	2.52	0.44
1:A:540:LEU:HG	1:A:601:VAL:HG11	1.98	0.44
1:B:85:CYS:SG	1:B:92:PHE:HB2	2.56	0.44
1:D:371:SER:HB3	1:D:374:ASP:HB2	1.99	0.44
1:D:395:VAL:HG13	1:D:473:ASP:HB2	1.98	0.44
1:D:402:GLU:HG3	1:D:450:TYR:OH	2.17	0.44
1:B:328:ALA:O	1:B:332:VAL:HG23	2.17	0.44
1:C:526:TRP:O	1:C:529:ILE:HG22	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:50:VAL:HG11	2:E:84:PRO:HA	1.99	0.44
1:B:97:PHE:HA	1:B:110:ARG:HD2	1.99	0.44
1:B:795:VAL:O	1:B:798:ILE:HG22	2.18	0.44
1:C:263:GLU:HG2	1:C:270:HIS:HB2	1.99	0.44
1:A:134:ASP:N	1:A:134:ASP:OD1	2.51	0.44
1:B:242:GLN:HE21	1:B:244:VAL:H	1.66	0.44
1:C:12:GLY:HA2	1:C:70:PHE:O	2.18	0.44
1:D:12:GLY:HA2	1:D:70:PHE:O	2.18	0.44
1:B:526:TRP:O	1:B:529:ILE:HG22	2.18	0.44
1:D:400:ILE:HG21	1:D:450:TYR:CZ	2.53	0.44
1:A:519:ASP:O	1:B:787:LEU:HD12	2.17	0.44
1:B:335:GLU:OE2	1:B:340:ASN:ND2	2.43	0.43
1:B:511:LYS:HA	1:B:512:PRO:HD2	1.83	0.43
1:B:356:ILE:O	1:B:356:ILE:HG13	2.18	0.43
1:C:85:CYS:SG	1:C:92:PHE:HB2	2.58	0.43
1:A:504:ILE:HD11	1:A:723:VAL:HG21	1.99	0.43
1:B:436:CYS:HA	2:F:58:MET:HE3	2.00	0.43
1:C:399:THR:OG1	1:C:406:VAL:HG21	2.19	0.43
1:A:611:ILE:HD13	1:B:795:VAL:HG21	2.00	0.43
1:C:625:THR:O	1:C:629:MET:HB2	2.19	0.43
1:C:633:ILE:HD13	1:C:633:ILE:HA	1.90	0.43
1:C:744:THR:HB	1:C:745:PRO:HD3	2.00	0.43
1:A:498:LEU:HD23	1:A:705:GLU:HG2	2.01	0.43
1:D:498:LEU:HB3	1:D:707:THR:HG23	2.01	0.43
2:F:21:LEU:HD21	2:F:32:VAL:HA	2.01	0.43
1:A:403:SER:HA	1:A:404:PRO:HA	1.74	0.43
1:B:97:PHE:HA	1:B:98:PRO:HD3	1.91	0.43
1:C:110:ARG:HA	1:C:111:PRO:HD3	1.88	0.43
1:C:457:THR:O	1:C:459:ILE:HG13	2.19	0.43
1:A:189:ILE:HG12	1:A:217:ILE:HB	2.01	0.43
1:A:464:VAL:O	1:A:468:VAL:HG23	2.19	0.43
2:E:8:ARG:CZ	2:E:12:CYS:HB2	2.48	0.43
1:A:795:VAL:HG21	1:D:611:ILE:HG21	2.00	0.42
1:C:243:ILE:HG23	1:C:244:VAL:HG23	2.00	0.42
1:C:504:ILE:HD13	1:C:504:ILE:H	1.83	0.42
3:D:901:KAI:HD2	3:D:901:KAI:HD12	1.77	0.42
1:C:809:VAL:HA	1:C:812:ILE:HG12	2.01	0.42
1:D:58:PHE:CE2	1:D:84:PHE:HB3	2.54	0.42
1:A:75:LYS:HB3	1:A:136:ASP:O	2.19	0.42
1:A:809:VAL:HA	1:A:812:ILE:HG12	2.01	0.42
1:B:73:TYR:CE2	1:B:94:THR:HG21	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:507:PRO:HG3	1:B:630:VAL:O	2.19	0.42
1:B:751:LEU:HA	5:B:903:FWF:H34	2.02	0.42
1:D:631:SER:HA	1:D:632:PRO:HD3	1.91	0.42
1:A:647:TYR:HB3	1:A:701:ALA:HB3	2.02	0.42
1:C:14:PHE:HA	1:C:15:PRO:HD3	1.92	0.42
1:B:22:TYR:CZ	1:B:43:ILE:HD13	2.55	0.42
1:C:137:ARG:HH22	1:C:192:CYS:C	2.23	0.42
2:E:38:GLU:HG2	2:E:38:GLU:H	1.69	0.42
1:D:89:HIS:HB3	1:D:322:GLY:HA3	2.02	0.42
1:D:625:THR:HG21	1:D:786:ALA:HB3	2.02	0.42
1:B:626:VAL:O	1:B:627:GLU:HG2	2.20	0.42
1:B:716:LYS:HA	1:B:718:CYS:H	1.85	0.42
5:B:903:FWF:N2	1:C:494:PRO:O	2.33	0.42
1:D:380:LEU:HG	1:D:381:THR:HG23	2.02	0.42
1:B:58:PHE:CE2	1:B:84:PHE:HB3	2.55	0.42
1:B:58:PHE:HE2	1:B:88:LEU:HD12	1.84	0.42
1:A:149:ALA:HA	1:A:154:TRP:HB2	2.02	0.42
1:B:647:TYR:HB3	1:B:701:ALA:HB3	2.02	0.42
1:C:395:VAL:HG13	1:C:473:ASP:HB2	2.01	0.42
1:A:126:TRP:CD2	1:A:187:ARG:HD3	2.55	0.42
1:B:243:ILE:HG23	1:B:244:VAL:HG23	2.02	0.41
1:D:44:ASP:OD1	1:D:61:GLN:NE2	2.53	0.41
1:D:526:TRP:O	1:D:529:ILE:HG22	2.20	0.41
1:D:19:ASP:HB3	1:D:267:PRO:HB2	2.01	0.41
1:A:475:ALA:HB3	1:A:735:ALA:HB3	2.02	0.41
1:B:716:LYS:HA	1:B:718:CYS:N	2.35	0.41
1:C:32:PHE:CZ	1:C:286:GLN:HB2	2.54	0.41
2:E:5:ASP:HB3	2:E:8:ARG:HB3	2.03	0.41
1:A:477:ALA:O	1:A:479:LEU:N	2.50	0.41
1:C:477:ALA:O	1:C:479:LEU:N	2.49	0.41
1:B:498:LEU:HB3	1:B:707:THR:HG23	2.03	0.41
1:C:795:VAL:O	1:C:798:ILE:HG22	2.20	0.41
2:F:20:CYS:SG	2:F:23:ARG:NH2	2.94	0.41
1:C:153:LYS:O	1:C:154:TRP:HD1	2.03	0.41
1:D:520:PRO:CG	1:D:620:LEU:HD23	2.51	0.41
2:E:16:ARG:NH1	2:E:42:THR:HG21	2.36	0.41
1:A:529:ILE:HG13	1:A:608:PHE:CZ	2.56	0.41
1:B:117:LEU:O	1:B:121:ILE:HG13	2.21	0.41
1:C:773:CYS:O	1:C:777:ASP:N	2.44	0.41
1:D:318:PRO:HG2	1:D:321:GLN:HG2	2.01	0.41
1:A:132:LEU:HD23	1:A:188:VAL:HG13	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:TYR:HA	1:A:176:LEU:HB3	2.03	0.41
1:A:399:THR:O	1:A:444:ILE:HA	2.21	0.41
1:A:494:PRO:O	5:D:903:FWF:N1	2.51	0.41
1:B:110:ARG:HA	1:B:111:PRO:HD3	1.82	0.41
1:D:97:PHE:HA	1:D:98:PRO:HD3	1.92	0.41
1:D:399:THR:O	1:D:444:ILE:HA	2.21	0.41
1:B:220:ASN:O	1:B:241:PHE:HB2	2.20	0.41
1:B:403:SER:HA	1:B:404:PRO:HA	1.75	0.41
2:E:80:SER:HA	2:F:5:ASP:OD2	2.21	0.41
1:A:220:ASN:O	1:A:241:PHE:HB2	2.21	0.40
1:A:526:TRP:O	1:A:529:ILE:HG22	2.21	0.40
1:C:618:ALA:HB1	1:D:621:ALA:HA	2.02	0.40
1:D:514:VAL:HG13	1:D:794:GLY:HA3	2.03	0.40
1:B:180:LEU:HD23	1:B:180:LEU:HA	1.90	0.40
3:A:901:KAI:HD2	3:A:901:KAI:HD12	1.68	0.40
1:D:450:TYR:CE2	3:D:901:KAI:HD1	2.57	0.40
1:B:499:GLY:O	1:B:706:SER:N	2.51	0.40
1:C:488:VAL:HG12	2:F:34:PHE:HD2	1.86	0.40
1:C:499:GLY:O	1:C:706:SER:N	2.54	0.40
1:C:536:VAL:HG21	1:C:605:TRP:CE3	2.56	0.40
1:D:397:VAL:HB	1:D:442:LEU:HD23	2.04	0.40
1:A:109:MET:O	1:A:349:ARG:NH1	2.53	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	730/814 (90%)	712 (98%)	18 (2%)	0	100	100
1	B	735/814 (90%)	714 (97%)	21 (3%)	0	100	100
1	C	740/814 (91%)	723 (98%)	17 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	730/814 (90%)	712 (98%)	18 (2%)	0	100	100
2	E	83/90 (92%)	77 (93%)	6 (7%)	0	100	100
2	F	83/90 (92%)	79 (95%)	4 (5%)	0	100	100
All	All	3101/3436 (90%)	3017 (97%)	84 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	563/695 (81%)	549 (98%)	14 (2%)	47	75
1	B	576/695 (83%)	567 (98%)	9 (2%)	62	83
1	C	542/695 (78%)	532 (98%)	10 (2%)	59	81
1	D	568/695 (82%)	558 (98%)	10 (2%)	59	81
2	E	73/76 (96%)	73 (100%)	0	100	100
2	F	73/76 (96%)	73 (100%)	0	100	100
All	All	2395/2932 (82%)	2352 (98%)	43 (2%)	59	81

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	73	TYR
1	A	75	LYS
1	A	89	HIS
1	A	210	HIS
1	A	255	ILE
1	A	271	THR
1	A	312	LEU
1	A	319	TRP
1	A	373	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	498	LEU
1	A	504	ILE
1	A	630	VAL
1	A	736	THR
1	B	73	TYR
1	B	192	CYS
1	B	271	THR
1	B	373	VAL
1	B	498	LEU
1	B	615	SER
1	B	619	ASN
1	B	772	GLU
1	B	787	LEU
1	C	46	LEU
1	C	73	TYR
1	C	166	ASN
1	C	192	CYS
1	C	271	THR
1	C	319	TRP
1	C	373	VAL
1	C	504	ILE
1	C	629	MET
1	C	638	ASP
1	D	44	ASP
1	D	46	LEU
1	D	73	TYR
1	D	271	THR
1	D	373	VAL
1	D	498	LEU
1	D	504	ILE
1	D	624	LEU
1	D	627	GLU
1	D	789	LEU

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

Mol	Chain	Res	Type
1	A	155	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

10 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	KAI	A	901	-	13,15,15	1.04	0	14,21,21	0.84	0
4	NAG	A	902	1	14,14,15	0.34	0	17,19,21	0.88	1 (5%)
4	NAG	D	902	1	14,14,15	0.37	0	17,19,21	0.71	1 (5%)
3	KAI	B	901	-	13,15,15	1.00	0	14,21,21	0.94	0
4	NAG	C	902	1	14,14,15	0.54	0	17,19,21	0.87	1 (5%)
4	NAG	B	902	1	14,14,15	0.18	0	17,19,21	0.66	0
5	FWF	D	903	-	31,33,33	1.48	2 (6%)	38,48,48	3.23	3 (7%)
3	KAI	C	901	-	13,15,15	1.07	0	14,21,21	0.92	0
3	KAI	D	901	-	13,15,15	1.06	0	14,21,21	0.88	0
5	FWF	B	903	-	31,33,33	1.47	2 (6%)	38,48,48	3.05	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
3	KAI	A	901	-	-	3/12/25/25	0/1/1/1
4	NAG	A	902	1	-	2/6/23/26	0/1/1/1
4	NAG	D	902	1	-	0/6/23/26	0/1/1/1
3	KAI	B	901	-	-	2/12/25/25	0/1/1/1
4	NAG	C	902	1	-	2/6/23/26	0/1/1/1
4	NAG	B	902	1	-	2/6/23/26	0/1/1/1
5	FWF	D	903	-	-	2/32/36/36	0/2/2/2
3	KAI	C	901	-	-	2/12/25/25	0/1/1/1
3	KAI	D	901	-	-	0/12/25/25	0/1/1/1
5	FWF	B	903	-	-	0/32/36/36	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	903	FWF	C19-S1	-5.72	1.66	1.78
5	B	903	FWF	C19-S1	-5.61	1.66	1.78
5	D	903	FWF	C20-S2	-5.59	1.66	1.78
5	B	903	FWF	C20-S2	-5.54	1.66	1.78

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	D	903	FWF	O2-S1-O1	-13.48	109.13	119.24
5	B	903	FWF	O4-S2-O3	-12.94	109.53	119.24
5	D	903	FWF	O4-S2-O3	-12.86	109.59	119.24
5	B	903	FWF	O2-S1-O1	-12.11	110.16	119.24
4	A	902	NAG	C1-O5-C5	3.07	116.35	112.19
4	C	902	NAG	C1-O5-C5	3.05	116.33	112.19
5	D	903	FWF	O1-S1-N2	2.49	110.76	107.76
5	B	903	FWF	O4-S2-N1	2.07	110.26	107.76
4	D	902	NAG	C1-O5-C5	2.07	115.00	112.19

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	902	NAG	C4-C5-C6-O6
4	B	902	NAG	C4-C5-C6-O6
4	C	902	NAG	O5-C5-C6-O6
4	B	902	NAG	O5-C5-C6-O6
4	A	902	NAG	C4-C5-C6-O6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
4	A	902	NAG	O5-C5-C6-O6
5	D	903	FWF	C22-C19-S1-O1
3	C	901	KAI	O-C-CA-N
3	C	901	KAI	OXT-C-CA-N
3	B	901	KAI	CB-CB1-CG1-OD1
3	B	901	KAI	CB-CB1-CG1-OD2
3	A	901	KAI	O-C-CA-CB
5	D	903	FWF	C21-C19-S1-O1
3	A	901	KAI	OXT-C-CA-CB
3	A	901	KAI	O-C-CA-N

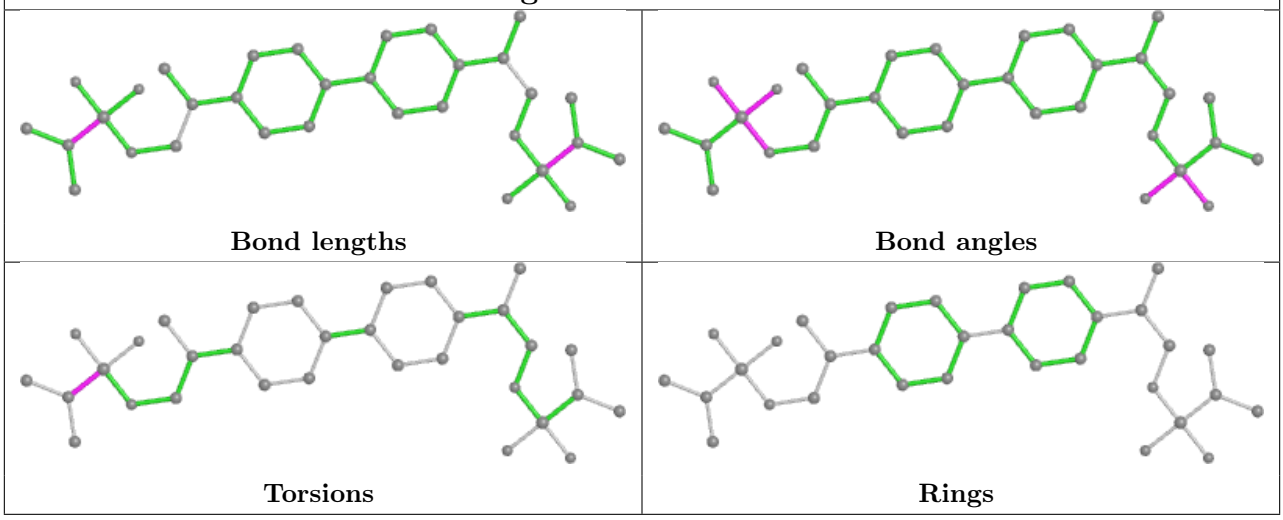
There are no ring outliers.

5 monomers are involved in 9 short contacts:

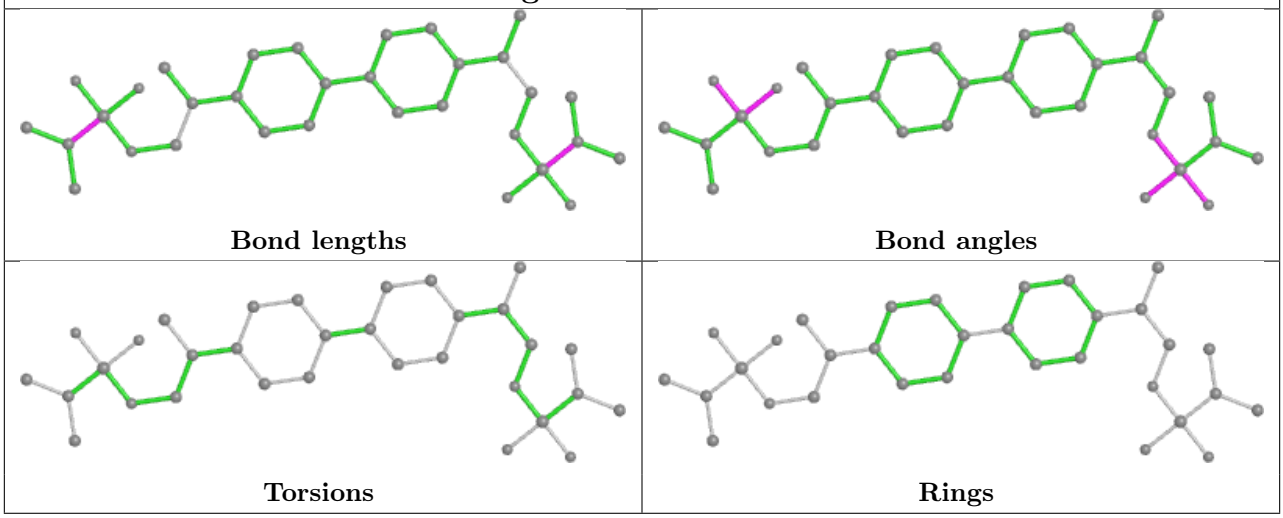
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	KAI	1	0
5	D	903	FWF	2	0
3	C	901	KAI	1	0
3	D	901	KAI	2	0
5	B	903	FWF	3	0

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 then 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 FWF D 903



Ligand FWF B 903



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 (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	738/814 (90%)	-0.18	10 (1%)	75	69	59, 118, 189, 245	0
1	B	743/814 (91%)	-0.16	9 (1%)	79	73	63, 117, 187, 223	0
1	C	746/814 (91%)	-0.22	10 (1%)	77	71	68, 122, 196, 277	0
1	D	738/814 (90%)	-0.18	9 (1%)	79	73	49, 116, 186, 247	0
2	E	85/90 (94%)	-0.25	0	100	100	57, 97, 142, 189	0
2	F	85/90 (94%)	-0.26	0	100	100	67, 105, 147, 190	0
All	All	3135/3436 (91%)	-0.19	38 (1%)	79	73	49, 117, 189, 277	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	630	VAL	5.2
1	D	516	SER	5.1
1	C	391	GLU	4.6
1	B	516	SER	4.1
1	A	631	SER	4.1
1	A	167	ASP	4.0
1	A	6	ASN	3.4
1	B	784	THR	3.3
1	B	314	ASN	3.3
1	D	626	VAL	3.2
1	C	784	THR	3.2
1	A	792	VAL	3.2
1	B	308	ALA	2.9
1	B	785	SER	2.9
1	B	620	LEU	2.8
1	C	12	GLY	2.8
1	D	792	VAL	2.8
1	B	791	ASN	2.8
1	D	515	PHE	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	392	GLN	2.7
1	D	627	GLU	2.7
1	A	210	HIS	2.6
1	D	603	GLY	2.6
1	C	782	GLU	2.5
1	B	460	TRP	2.5
1	C	510	SER	2.5
1	C	783	LYS	2.4
1	D	604	VAL	2.3
1	A	305	ARG	2.3
1	A	137	ARG	2.1
1	C	209	LYS	2.1
1	C	629	MET	2.1
1	C	129	PHE	2.1
1	A	306	GLY	2.1
1	D	606	TRP	2.0
1	C	212	LYS	2.0
1	D	512	PRO	2.0
1	A	516	SER	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NAG	D	902	14/15	0.80	0.28	153,205,220,238	0
4	NAG	B	902	14/15	0.85	0.22	159,180,190,213	0
4	NAG	A	902	14/15	0.92	0.25	65,115,137,155	0
5	FWF	D	903	32/32	0.92	0.39	5,59,323,419	0

Continued on next page...

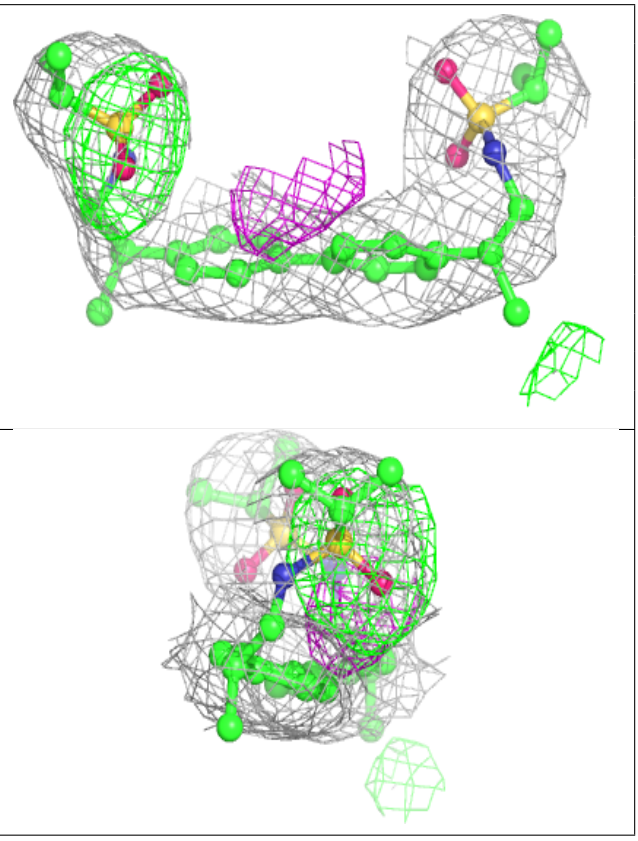
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	KAI	C	901	15/15	0.93	0.26	40,90,148,157	0
3	KAI	B	901	15/15	0.94	0.21	52,83,141,158	0
4	NAG	C	902	14/15	0.94	0.23	49,105,142,142	0
5	FWF	B	903	32/32	0.95	0.43	26,98,126,141	0
3	KAI	A	901	15/15	0.96	0.31	39,96,140,147	0
3	KAI	D	901	15/15	0.96	0.30	7,52,122,166	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.

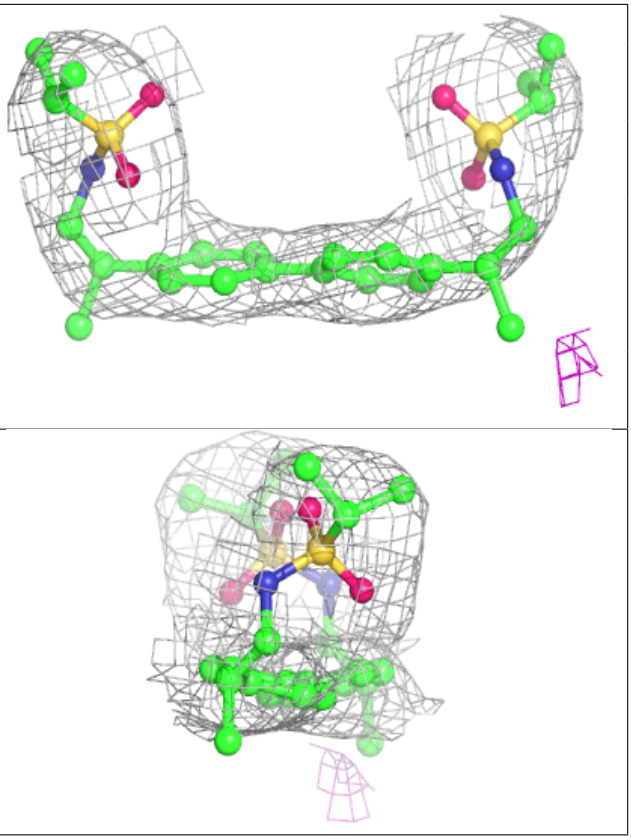
Electron density around FWF D 903:

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 FWF B 903:

$2mF_o$-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.