



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 07:53 AM UTC

PDB ID : 1SFC / pdb_00001sfc
Title : NEURONAL SYNAPTIC FUSION COMPLEX
Authors : Sutton, R.B.; Brunger, A.T.
Deposited on : 1998-08-24
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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)
Xtrriage (Phenix) : 2.0
EDS : 3.0
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

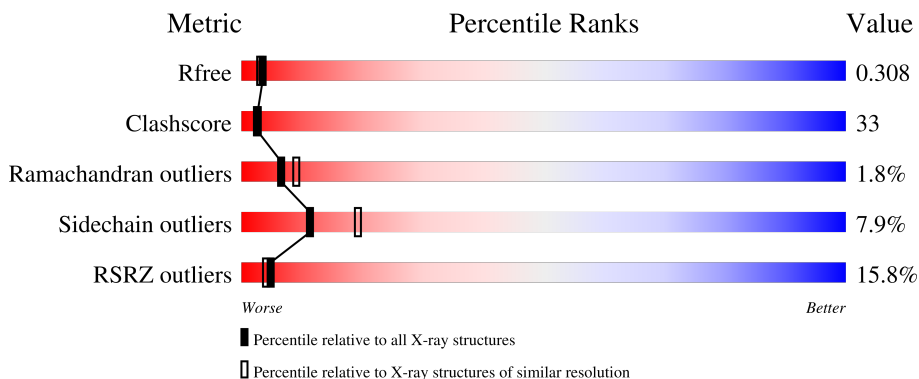
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



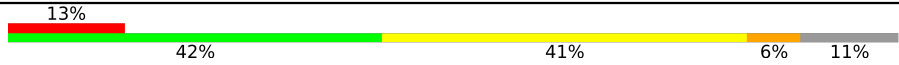
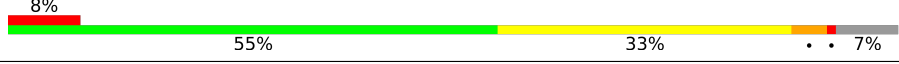
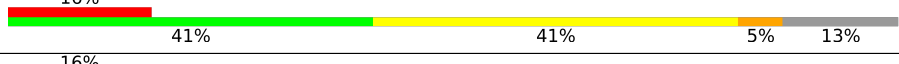

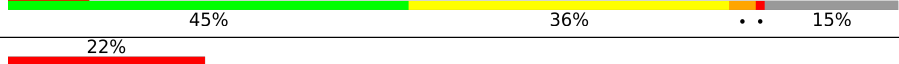
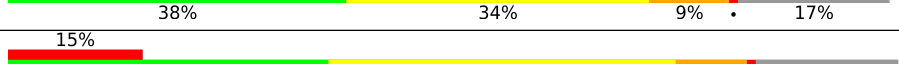
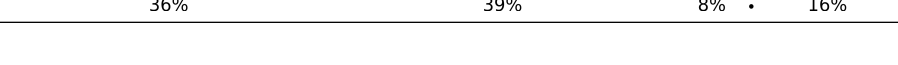
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 $\leq 5\%$. 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	96	
1	E	96	
1	I	96	
2	B	83	
2	F	83	

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Mol	Chain	Length	Quality of chain
2	J	83	
3	C	83	
3	G	83	
3	K	83	
4	D	87	
4	H	87	
4	L	87	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7105 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 PROTEIN (SYNAPTOBREVIN 2).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	69	564	347	106	110	1	0	0	0
1	E	69	567	349	106	109	3	0	0	0
1	I	69	565	349	106	109	1	0	0	0

- Molecule 2 is a protein called PROTEIN (SYNTAXIN 1A).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	72	586	362	100	119	5	0	0	0
2	F	73	594	368	102	119	5	0	0	0
2	J	74	600	371	102	122	5	20	0	0

- Molecule 3 is a protein called PROTEIN (SNAP-25B).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	77	625	371	115	133	6	0	0	0
3	G	72	581	345	106	125	5	9	0	0
3	K	73	589	351	107	126	5	9	0	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	58	ASP	GLU	SEE REMARK 999	UNP P60881
C	60	VAL	ILE	SEE REMARK 999	UNP P60881

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Chain	Residue	Modelled	Actual	Comment	Reference
C	65	ASN	ASP	SEE REMARK 999	UNP P60881
C	66	HIS	GLN	SEE REMARK 999	UNP P60881
C	69	GLN	LYS	SEE REMARK 999	UNP P60881
C	79	LYS	THR	SEE REMARK 999	UNP P60881
G	58	ASP	GLU	SEE REMARK 999	UNP P60881
G	60	VAL	ILE	SEE REMARK 999	UNP P60881
G	65	ASN	ASP	SEE REMARK 999	UNP P60881
G	66	HIS	GLN	SEE REMARK 999	UNP P60881
G	69	GLN	LYS	SEE REMARK 999	UNP P60881
G	79	LYS	THR	SEE REMARK 999	UNP P60881
K	58	ASP	GLU	SEE REMARK 999	UNP P60881
K	60	VAL	ILE	SEE REMARK 999	UNP P60881
K	65	ASN	ASP	SEE REMARK 999	UNP P60881
K	66	HIS	GLN	SEE REMARK 999	UNP P60881
K	69	GLN	LYS	SEE REMARK 999	UNP P60881
K	79	LYS	THR	SEE REMARK 999	UNP P60881

- Molecule 4 is a protein called PROTEIN (SNAP-25B).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	74	Total	C	N	O	S	11	0	0
			586	346	115	120	5			
4	H	72	Total	C	N	O	S	0	0	0
			578	342	113	118	5			
4	L	73	Total	C	N	O	S	7	0	0
			582	344	114	119	5			

- Molecule 5 is STRONTIUM ION (CCD ID: SR) (formula: Sr).

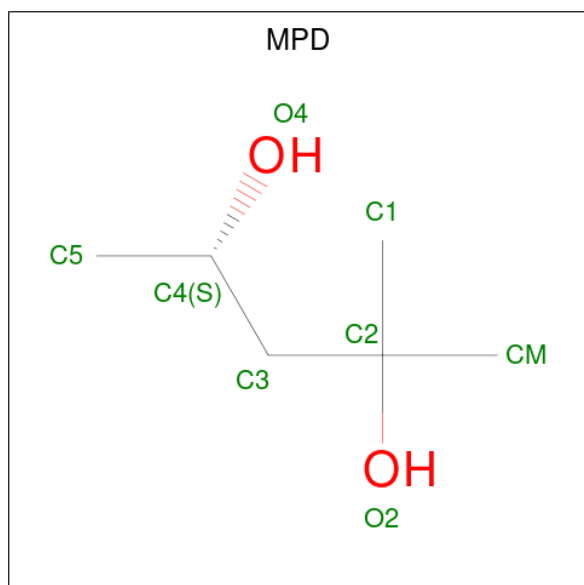
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Sr	0	0
			2	2		
5	B	1	Total	Sr	0	0
			1	1		
5	C	2	Total	Sr	0	0
			2	2		
5	D	1	Total	Sr	0	0
			1	1		
5	F	2	Total	Sr	0	0
			2	2		
5	G	3	Total	Sr	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	1	Total	Sr	0	0
			1	1		
5	K	1	Total	Sr	0	0
			1	1		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			8	6	2		
6	A	1	Total	C	O	0	0
			8	6	2		
6	F	1	Total	C	O	0	0
			8	6	2		
6	G	1	Total	C	O	0	0
			8	6	2		
6	I	1	Total	C	O	0	0
			8	6	2		
6	K	1	Total	C	O	0	0
			8	6	2		
6	K	1	Total	C	O	0	0
			8	6	2		

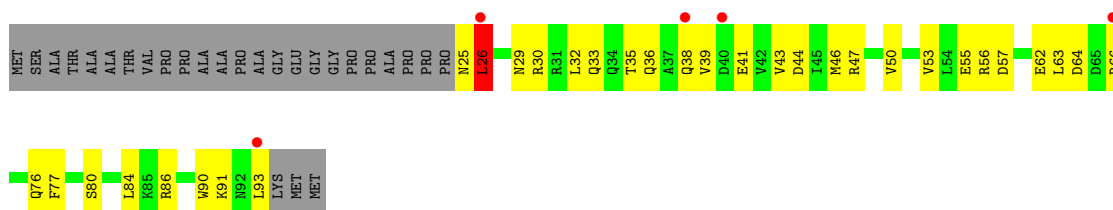
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	2	Total O 2 2	0	0
7	B	2	Total O 2 2	0	0
7	C	1	Total O 1 1	0	0
7	D	1	Total O 1 1	0	0
7	F	5	Total O 5 5	0	0
7	G	1	Total O 1 1	0	0
7	I	3	Total O 3 3	0	0
7	J	3	Total O 3 3	0	0
7	K	1	Total O 1 1	0	0

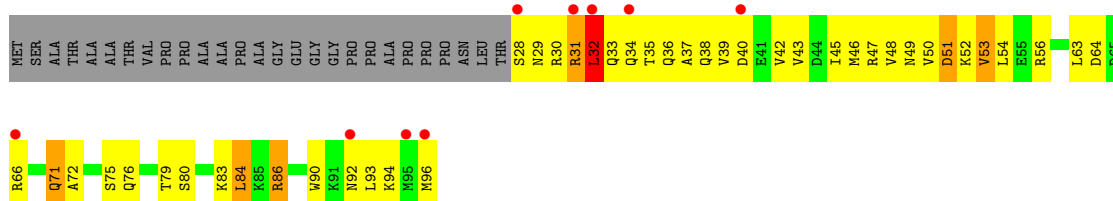
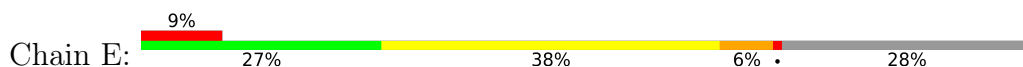
3 Residue-property plots [i](#)

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.

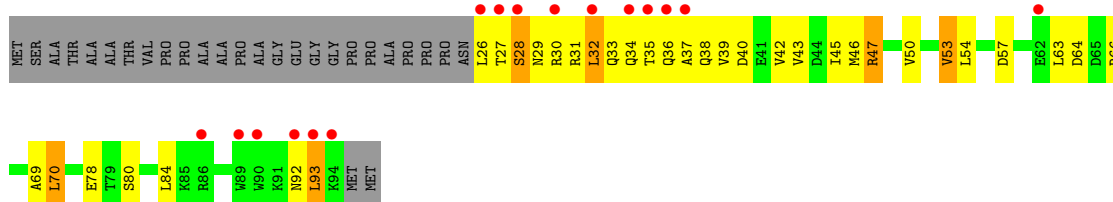
- Molecule 1: PROTEIN (SYNAPTOBREVIN 2)



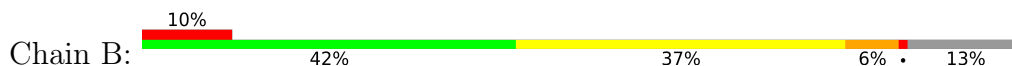
- Molecule 1: PROTEIN (SYNAPTOBREVIN 2)



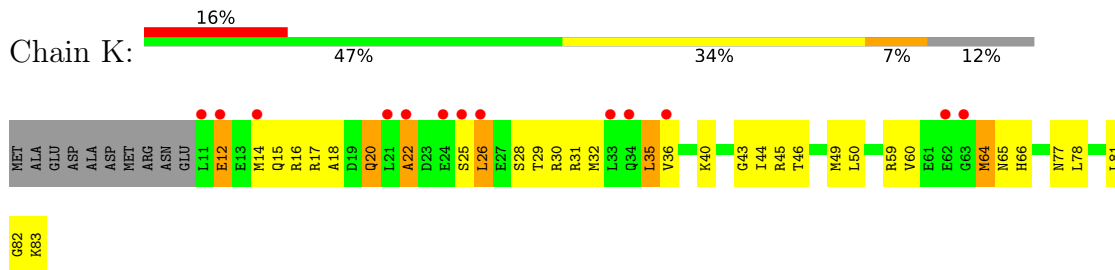
- Molecule 1: PROTEIN (SYNAPTOBREVIN 2)



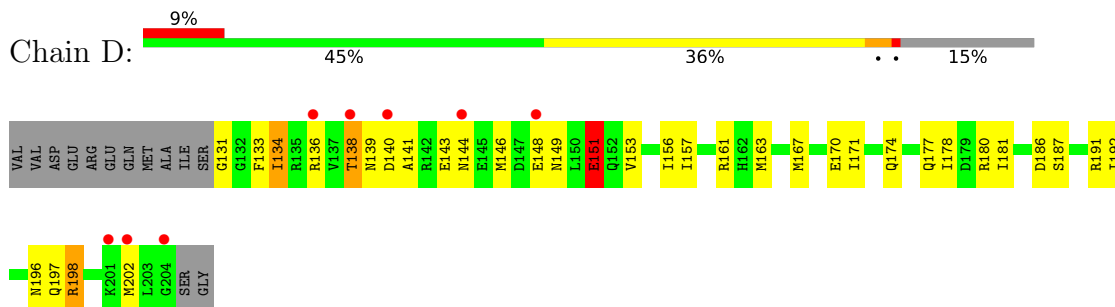
- Molecule 2: PROTEIN (SYNTAXIN 1A)



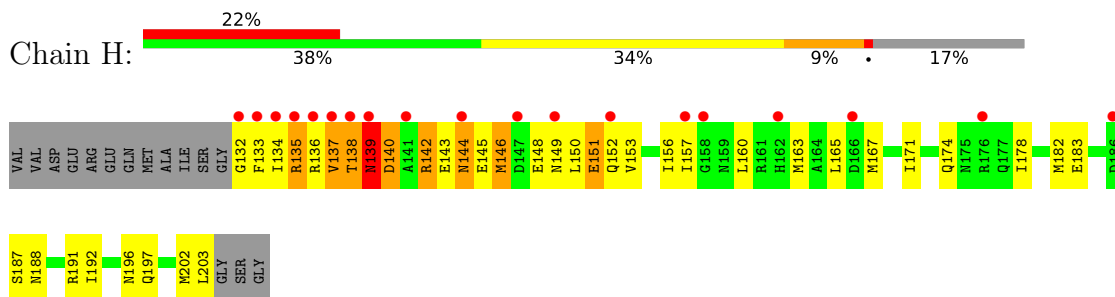
- Molecule 3: PROTEIN (SNAP-25B)



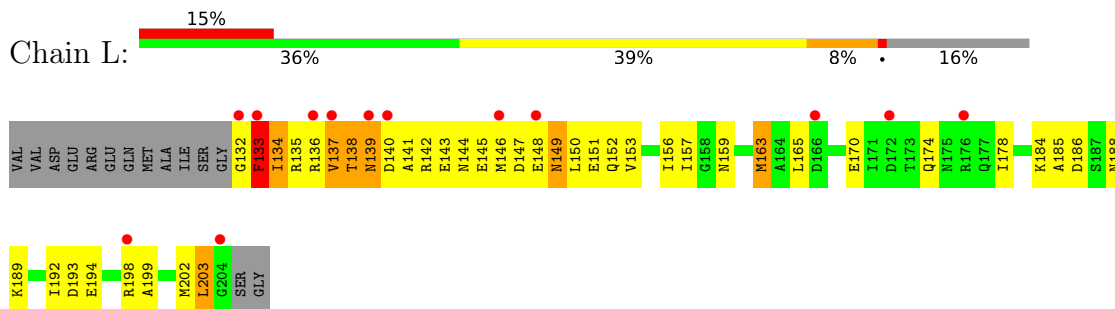
- Molecule 4: PROTEIN (SNAP-25B)



- Molecule 4: PROTEIN (SNAP-25B)



- Molecule 4: PROTEIN (SNAP-25B)



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	98.72Å 111.07Å 198.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.8 (50.00-2.40) 98.6 (50.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.47 (at 2.39Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.265 , 0.303 0.271 , 0.308	Depositor DCC
R_{free} test set	8052 reflections (9.77%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtrriage
Anisotropy	0.336	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.21$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7105	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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: SR, MPD

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	0.49	0/569	1.02	3/765 (0.4%)
1	E	0.44	0/572	0.92	1/764 (0.1%)
1	I	0.51	0/570	1.00	2/765 (0.3%)
2	B	0.55	0/592	1.08	3/792 (0.4%)
2	F	0.51	0/600	1.00	2/802 (0.2%)
2	J	0.52	0/606	1.01	3/811 (0.4%)
3	C	0.53	0/625	0.92	1/828 (0.1%)
3	G	0.52	0/581	1.14	4/770 (0.5%)
3	K	0.51	0/589	1.03	5/781 (0.6%)
4	D	0.55	1/587 (0.2%)	1.00	2/782 (0.3%)
4	H	0.62	1/579 (0.2%)	1.06	4/772 (0.5%)
4	L	0.61	1/583 (0.2%)	1.06	4/777 (0.5%)
All	All	0.53	3/7053 (0.0%)	1.02	34/9409 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	163	MET	SD-CE	-6.59	1.63	1.79
4	H	146	MET	SD-CE	-5.46	1.66	1.79
4	D	131	GLY	N-CA	5.34	1.54	1.45

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	38	GLU	N-CA-C	-11.96	99.16	113.88
2	B	190	GLN	N-CA-C	-11.72	98.69	113.23
3	G	14	MET	N-CA-C	-8.96	99.24	111.87
3	G	16	ARG	N-CA-C	-8.69	102.58	113.01
2	F	232	ARG	N-CA-C	-7.90	101.97	111.69

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	564	0	565	29	0
1	E	567	0	573	71	0
1	I	565	0	571	60	0
2	B	586	0	575	49	0
2	F	594	0	588	68	0
2	J	600	0	591	56	0
3	C	625	0	614	44	0
3	G	581	0	569	69	0
3	K	589	0	580	57	0
4	D	586	0	573	33	0
4	H	578	0	567	86	0
4	L	582	0	570	74	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
5	D	1	0	0	0	0
5	F	2	0	0	0	0
5	G	3	0	0	0	0
5	J	1	0	0	0	0
5	K	1	0	0	0	0
6	A	16	0	28	0	0
6	F	8	0	14	0	0
6	G	8	0	14	1	0
6	I	8	0	14	2	0
6	K	16	0	28	1	0
7	A	2	0	0	0	0
7	B	2	0	0	2	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
7	F	5	0	0	0	0
7	G	1	0	0	0	0
7	I	3	0	0	0	0
7	J	3	0	0	0	0
7	K	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7105	0	7034	469	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 33.

The worst 5 of 469 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)
3:G:12:GLU:HG2	3:G:13:GLU:H	1.20	1.01
3:C:71:MET:HE3	4:D:192:ILE:HG12	1.44	0.98
1:E:53:VAL:HG21	2:F:219:MET:SD	2.03	0.98
2:B:221:MET:HE1	1:I:93:LEU:HD21	1.47	0.96
1:I:31:ARG:HG2	4:L:132:GLY:HA2	1.47	0.95

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	67/96 (70%)	64 (96%)	2 (3%)	1 (2%)	8	12
1	E	67/96 (70%)	63 (94%)	3 (4%)	1 (2%)	8	12
1	I	67/96 (70%)	63 (94%)	3 (4%)	1 (2%)	8	12
2	B	70/83 (84%)	68 (97%)	2 (3%)	0	100	100
2	F	71/83 (86%)	67 (94%)	2 (3%)	2 (3%)	4	3
2	J	72/83 (87%)	68 (94%)	4 (6%)	0	100	100
3	C	75/83 (90%)	73 (97%)	2 (3%)	0	100	100
3	G	70/83 (84%)	67 (96%)	3 (4%)	0	100	100
3	K	71/83 (86%)	68 (96%)	3 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	D	72/87 (83%)	63 (88%)	7 (10%)	2 (3%)	4	3
4	H	70/87 (80%)	61 (87%)	6 (9%)	3 (4%)	2	1
4	L	71/87 (82%)	62 (87%)	4 (6%)	5 (7%)	1	0
All	All	843/1047 (80%)	787 (93%)	41 (5%)	15 (2%)	6	9

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
4	H	139	ASN
1	I	93	LEU
4	L	138	THR
4	D	139	ASN

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	61/78 (78%)	58 (95%)	3 (5%)	22	39
1	E	61/78 (78%)	55 (90%)	6 (10%)	7	12
1	I	61/78 (78%)	56 (92%)	5 (8%)	10	18
2	B	66/75 (88%)	56 (85%)	10 (15%)	3	3
2	F	66/75 (88%)	62 (94%)	4 (6%)	17	30
2	J	68/75 (91%)	65 (96%)	3 (4%)	25	43
3	C	69/73 (94%)	60 (87%)	9 (13%)	4	5
3	G	64/73 (88%)	57 (89%)	7 (11%)	6	9
3	K	65/73 (89%)	62 (95%)	3 (5%)	24	41
4	D	63/74 (85%)	60 (95%)	3 (5%)	23	40
4	H	63/74 (85%)	57 (90%)	6 (10%)	8	13
4	L	63/74 (85%)	61 (97%)	2 (3%)	34	56
All	All	770/900 (86%)	709 (92%)	61 (8%)	11	19

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	71	GLN
2	J	259	SER
3	G	13	GLU
2	J	251	THR
4	L	133	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	36	GLN
3	K	20	GLN
1	I	49	ASN
2	J	213	HIS
3	K	66	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 13 are monoatomic - leaving 7 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
6	MPD	K	929	-	7,7,7	1.25	1 (14%)	9,10,10	2.90	4 (44%)
6	MPD	K	931	-	7,7,7	1.18	1 (14%)	9,10,10	2.79	3 (33%)
6	MPD	F	926	-	7,7,7	1.10	1 (14%)	9,10,10	2.94	3 (33%)
6	MPD	I	930	-	7,7,7	1.20	1 (14%)	9,10,10	2.88	4 (44%)
6	MPD	A	927	-	7,7,7	1.18	1 (14%)	9,10,10	2.71	3 (33%)
6	MPD	A	928	-	7,7,7	1.20	1 (14%)	9,10,10	2.80	4 (44%)
6	MPD	G	932	-	7,7,7	1.13	1 (14%)	9,10,10	2.78	3 (33%)

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
6	MPD	K	929	-	-	1/5/5/5	-
6	MPD	K	931	-	-	0/5/5/5	-
6	MPD	F	926	-	-	0/5/5/5	-
6	MPD	I	930	-	-	0/5/5/5	-
6	MPD	A	927	-	-	0/5/5/5	-
6	MPD	A	928	-	-	0/5/5/5	-
6	MPD	G	932	-	-	0/5/5/5	-

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	K	929	MPD	O2-C2	2.68	1.51	1.44
6	K	931	MPD	O2-C2	2.65	1.51	1.44
6	A	927	MPD	O2-C2	2.64	1.51	1.44
6	A	928	MPD	O2-C2	2.54	1.50	1.44
6	I	930	MPD	O2-C2	2.54	1.50	1.44

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	926	MPD	CM-C2-C1	7.06	126.43	110.63
6	K	929	MPD	CM-C2-C1	7.02	126.35	110.63
6	I	930	MPD	CM-C2-C1	6.93	126.16	110.63
6	A	928	MPD	CM-C2-C1	6.73	125.71	110.63
6	K	931	MPD	CM-C2-C1	6.72	125.67	110.63

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	K	929	MPD	C1-C2-C3-C4

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	K	931	MPD	1	0
6	I	930	MPD	2	0
6	G	932	MPD	1	0

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	69/96 (71%)	0.78	5 (7%) 21 18	28, 39, 57, 64	0
1	E	69/96 (71%)	0.88	9 (13%) 7 5	23, 41, 73, 84	0
1	I	69/96 (71%)	1.04	16 (23%) 2 1	19, 39, 76, 80	0
2	B	72/83 (86%)	0.71	8 (11%) 10 7	20, 36, 60, 74	0
2	F	73/83 (87%)	1.04	14 (19%) 3 2	22, 41, 82, 89	0
2	J	71/83 (85%)	0.90	11 (15%) 5 4	20, 38, 76, 82	0
3	C	77/83 (92%)	0.70	7 (9%) 15 12	22, 39, 73, 79	0
3	G	72/83 (86%)	1.03	13 (18%) 3 3	0, 42, 91, 94	1 (1%)
3	K	72/83 (86%)	0.93	13 (18%) 3 3	20, 42, 70, 74	0
4	D	73/87 (83%)	0.80	8 (10%) 10 8	24, 40, 63, 70	0
4	H	72/87 (82%)	1.33	19 (26%) 1 1	26, 45, 83, 86	0
4	L	72/87 (82%)	0.99	13 (18%) 3 3	22, 44, 70, 72	0
All	All	861/1047 (82%)	0.92	136 (15%) 5 4	0, 41, 76, 94	1 (0%)

The worst 5 of 136 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	261	ALA	6.5
4	H	137	VAL	6.1
4	H	132	GLY	4.9
1	I	27	THR	4.9
3	G	82	GLY	4.6

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(Å ²)	Q<0.9
6	MPD	K	929	8/8	0.34	0.28	79,80,80,81	0
6	MPD	I	930	8/8	0.73	0.24	65,69,70,72	0
6	MPD	G	932	8/8	0.75	0.24	70,71,72,72	0
5	SR	C	295	1/1	0.77	0.14	87,87,87,87	1
6	MPD	F	926	8/8	0.83	0.20	44,46,50,52	0
5	SR	K	293	1/1	0.83	0.10	96,96,96,96	1
5	SR	F	297	1/1	0.84	0.09	95,95,95,95	0
5	SR	C	296	1/1	0.85	0.09	96,96,96,96	1
6	MPD	A	927	8/8	0.86	0.18	74,75,75,75	0
6	MPD	K	931	8/8	0.87	0.14	70,71,71,72	0
6	MPD	A	928	8/8	0.89	0.19	51,54,55,56	0
5	SR	B	294	1/1	0.90	0.07	96,96,96,96	1
5	SR	G	299	1/1	0.90	0.11	92,92,92,92	0
5	SR	G	300	1/1	0.91	0.10	81,81,81,81	0
5	SR	J	292	1/1	0.92	0.09	96,96,96,96	0
5	SR	G	298	1/1	0.95	0.08	63,63,63,63	0
5	SR	D	288	1/1	0.96	0.12	58,58,58,58	0
5	SR	F	291	1/1	0.98	0.07	63,63,63,63	0
5	SR	A	290	1/1	0.98	0.06	50,50,50,50	0
5	SR	A	289	1/1	1.00	0.02	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.