



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:18 pm GMT

PDB ID : 4BUO
Title : High Resolution Structure of Thermostable Agonist-bound Neurotensin Receptor 1 Mutant without Lysozyme Fusion
Authors : Egloff, P.; Hillenbrand, M.; Schlinkmann, K.M.; Batyuk, A.; Mittl, P.; Plueckthun, A.
Deposited on : 2013-06-21
Resolution : 2.75 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

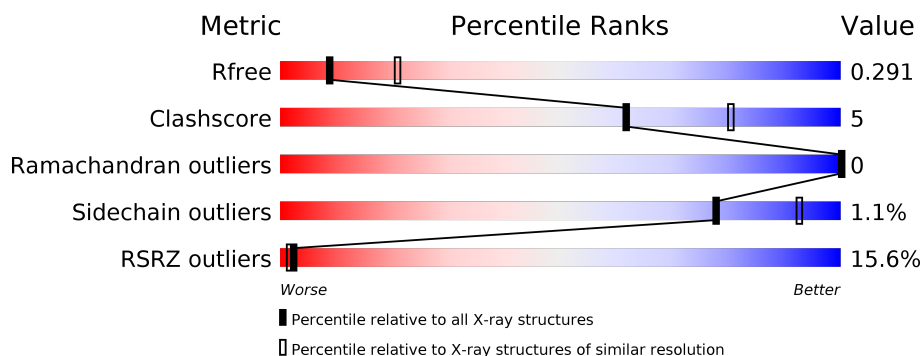
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	100719	3666 (2.80-2.72)
Clashscore	112137	4174 (2.80-2.72)
Ramachandran outliers	110173	4103 (2.80-2.72)
Sidechain outliers	110143	4106 (2.80-2.72)
RSRZ outliers	101464	3697 (2.80-2.72)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	335	
1	B	335	
2	C	10	
2	D	10	

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

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GLY	A	1385	-	-	-	X
3	GLY	A	1387	-	-	-	X
3	GLY	A	1392	-	-	-	X
3	GLY	B	1387	-	-	-	X
3	GLY	B	1388	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5034 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 NEUROTENSIN RECEPTOR TYPE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2400	1589	387	409	15			
1	B	314	Total	C	N	O	S	0	0	0
			2463	1628	394	424	17			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	PRO	-	EXPRESSION TAG	UNP P20789
A	48	GLY	-	EXPRESSION TAG	UNP P20789
A	49	SER	-	EXPRESSION TAG	UNP P20789
A	391	THR	-	EXPRESSION TAG	UNP P20789
A	392	ARG	-	EXPRESSION TAG	UNP P20789
A	393	GLU	-	EXPRESSION TAG	UNP P20789
A	394	LEU	-	EXPRESSION TAG	UNP P20789
A	395	GLU	-	EXPRESSION TAG	UNP P20789
A	396	VAL	-	EXPRESSION TAG	UNP P20789
A	397	LEU	-	EXPRESSION TAG	UNP P20789
A	398	PHE	-	EXPRESSION TAG	UNP P20789
A	399	GLN	-	EXPRESSION TAG	UNP P20789
A	86	LEU	ALA	ENGINEERED MUTATION	UNP P20789
A	103	ASP	HIS	ENGINEERED MUTATION	UNP P20789
A	105	TYR	HIS	ENGINEERED MUTATION	UNP P20789
A	161	VAL	ALA	ENGINEERED MUTATION	UNP P20789
A	167	LEU	ARG	ENGINEERED MUTATION	UNP P20789
A	213	LEU	ARG	ENGINEERED MUTATION	UNP P20789
A	234	LEU	VAL	ENGINEERED MUTATION	UNP P20789
A	253	ALA	ILE	ENGINEERED MUTATION	UNP P20789
A	305	ARG	HIS	ENGINEERED MUTATION	UNP P20789
A	358	VAL	PHE	ENGINEERED MUTATION	UNP P20789
A	362	ALA	SER	ENGINEERED MUTATION	UNP P20789
B	47	PRO	-	EXPRESSION TAG	UNP P20789
B	48	GLY	-	EXPRESSION TAG	UNP P20789

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Chain	Residue	Modelled	Actual	Comment	Reference
B	49	SER	-	EXPRESSION TAG	UNP P20789
B	391	THR	-	EXPRESSION TAG	UNP P20789
B	392	ARG	-	EXPRESSION TAG	UNP P20789
B	393	GLU	-	EXPRESSION TAG	UNP P20789
B	394	LEU	-	EXPRESSION TAG	UNP P20789
B	395	GLU	-	EXPRESSION TAG	UNP P20789
B	396	VAL	-	EXPRESSION TAG	UNP P20789
B	397	LEU	-	EXPRESSION TAG	UNP P20789
B	398	PHE	-	EXPRESSION TAG	UNP P20789
B	399	GLN	-	EXPRESSION TAG	UNP P20789
B	86	LEU	ALA	ENGINEERED MUTATION	UNP P20789
B	103	ASP	HIS	ENGINEERED MUTATION	UNP P20789
B	105	TYR	HIS	ENGINEERED MUTATION	UNP P20789
B	161	VAL	ALA	ENGINEERED MUTATION	UNP P20789
B	167	LEU	ARG	ENGINEERED MUTATION	UNP P20789
B	213	LEU	ARG	ENGINEERED MUTATION	UNP P20789
B	234	LEU	VAL	ENGINEERED MUTATION	UNP P20789
B	253	ALA	ILE	ENGINEERED MUTATION	UNP P20789
B	305	ARG	HIS	ENGINEERED MUTATION	UNP P20789
B	358	VAL	PHE	ENGINEERED MUTATION	UNP P20789
B	362	ALA	SER	ENGINEERED MUTATION	UNP P20789

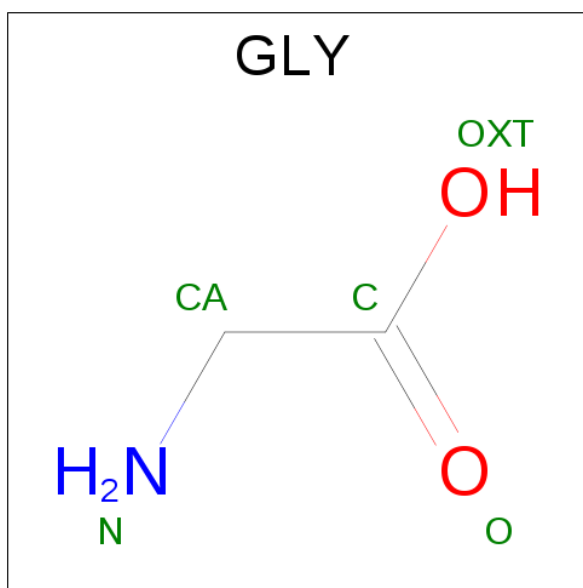
- Molecule 2 is a protein called NEUROTENSIN/NEUROMEDIN N.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			58	38	12	8			
2	D	6	Total	C	N	O	0	0	0
			58	38	12	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	GLY	-	SEE REMARK 999	UNP P20068
C	5	PRO	-	SEE REMARK 999	UNP P20068
C	6	GLY	-	SEE REMARK 999	UNP P20068
C	7	GLY	-	SEE REMARK 999	UNP P20068
D	4	GLY	-	SEE REMARK 999	UNP P20068
D	5	PRO	-	SEE REMARK 999	UNP P20068
D	6	GLY	-	SEE REMARK 999	UNP P20068
D	7	GLY	-	SEE REMARK 999	UNP P20068

- Molecule 3 is GLYCINE (three-letter code: GLY) (formula: $C_2H_5NO_2$).

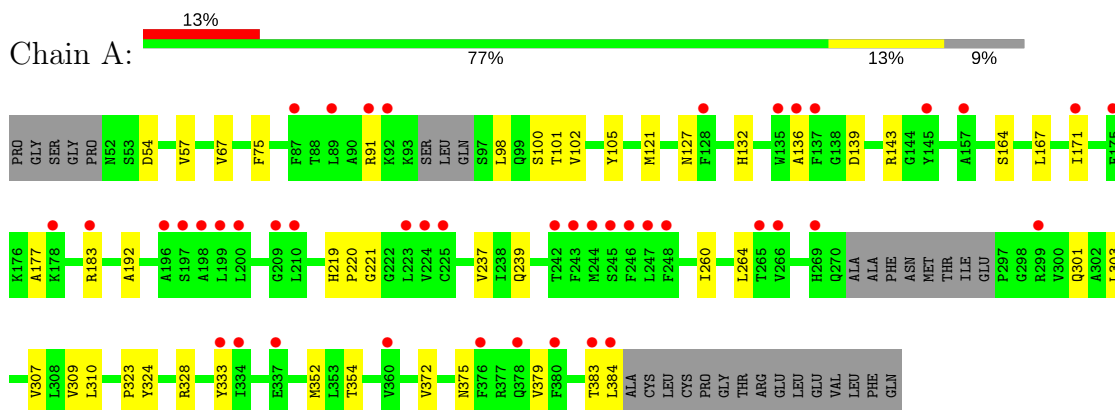


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	A	1	Total	C	N	O	0	0
			5	2	1	2		
3	B	1	Total	C	N	O	0	0
			5	2	1	2		
3	B	1	Total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

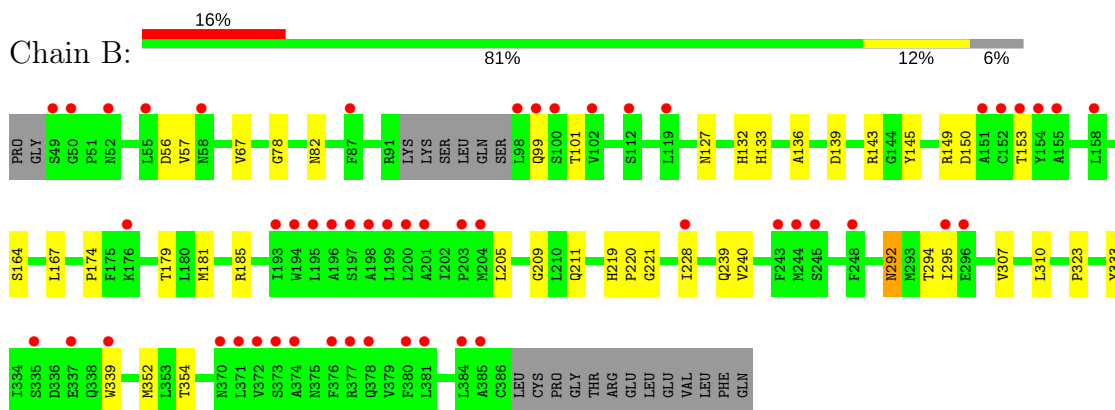
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

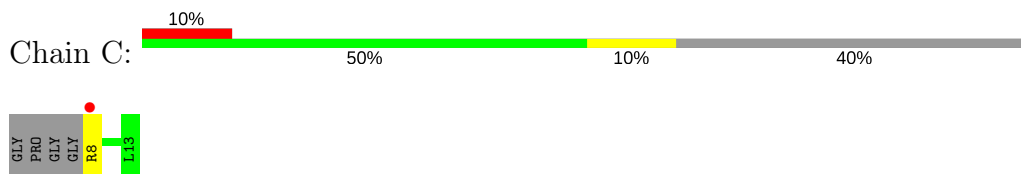
• Molecule 1: NEUROTENSIN RECEPTOR TYPE 1



• Molecule 1: NEUROTENSIN RECEPTOR TYPE 1



• Molecule 2: NEUROTENSIN/NEUROMEDIN N



• Molecule 2: NEUROTENSIN/NEUROMEDIN N





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.31Å 89.40Å 212.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.82 – 2.75 46.45 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (19.82-2.75) 98.8 (46.45-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.77Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.248 , 0.273 0.256 , 0.291	Depositor DCC
R_{free} test set	1587 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	110.0	Xtriage
Anisotropy	0.289	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5034	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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](#)

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	0/2459	0.39	0/3355
1	B	0.23	0/2525	0.38	0/3451
2	C	0.18	0/59	0.38	0/77
2	D	0.18	0/59	0.38	0/77
All	All	0.23	0/5102	0.39	0/6960

There are no bond length outliers.

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	2400	0	2465	27	0
1	B	2463	0	2512	26	0
2	C	58	0	63	1	0
2	D	58	0	63	2	0
3	A	40	0	16	0	0
3	B	15	0	6	1	0
All	All	5034	0	5125	52	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 (52) 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:91:ARG:HH21	1:A:379:VAL:HG12	1.50	0.77
1:A:127:ASN:HD21	1:A:136:ALA:H	1.39	0.70
1:A:239:GLN:NE2	1:A:333:TYR:OH	2.25	0.69
1:B:56:ASP:O	2:D:8:ARG:NH2	2.26	0.69
1:B:239:GLN:NE2	1:B:333:TYR:OH	2.26	0.68
1:A:219:HIS:HD2	1:A:221:GLY:H	1.41	0.67
1:B:127:ASN:HD21	1:B:136:ALA:H	1.43	0.65
1:B:219:HIS:HD2	1:B:221:GLY:H	1.45	0.62
1:B:139:ASP:OD2	1:B:143:ARG:NH1	2.33	0.60
1:A:264:LEU:HD22	1:A:307:VAL:HG21	1.85	0.59
1:A:323:PRO:HB2	1:A:354:THR:HG22	1.88	0.56
1:A:98:LEU:HD21	1:A:177:ALA:HB1	1.90	0.54
1:A:167:LEU:HD23	1:A:260:ILE:HD13	1.91	0.53
1:A:192:ALA:HA	1:B:240:VAL:HG22	1.91	0.53
1:A:100:SER:HB3	1:A:183:ARG:HH21	1.75	0.52
1:A:219:HIS:CD2	1:A:221:GLY:H	2.26	0.52
1:B:78:GLY:O	1:B:82:ASN:ND2	2.40	0.51
1:B:294:THR:OG1	1:B:295:ILE:N	2.37	0.51
1:B:174:PRO:HB3	1:B:295:ILE:HD13	1.93	0.50
1:A:54:ASP:O	2:C:8:ARG:NH1	2.45	0.50
1:B:219:HIS:CD2	1:B:221:GLY:H	2.29	0.49
1:A:67:VAL:HG21	1:A:352:MET:HB3	1.94	0.49
1:A:171:ILE:HG22	1:A:303:LEU:HG	1.95	0.49
1:A:375:ASN:O	1:A:379:VAL:HG23	2.14	0.48
1:B:211:GLN:HG3	1:B:228:ILE:HG22	1.95	0.48
1:A:139:ASP:OD2	1:A:143:ARG:NH1	2.47	0.48
1:A:98:LEU:O	1:A:101:THR:OG1	2.20	0.48
1:B:209:GLY:HA3	1:B:228:ILE:HD13	1.97	0.47
1:B:145:TYR:CZ	1:B:149:ARG:HD2	2.50	0.46
1:B:57:VAL:HB	1:B:132:HIS:CD2	2.50	0.46
1:A:309:VAL:CG2	1:A:372:VAL:HG11	2.46	0.45
1:B:99:GLN:NE2	3:B:1388:GLY:O	2.50	0.45
1:B:219:HIS:HA	1:B:220:PRO:HD3	1.85	0.44
1:B:181:MET:HB3	1:B:185:ARG:HB2	2.00	0.44
1:B:143:ARG:NE	1:B:205:LEU:O	2.43	0.44
1:B:101:THR:OG1	1:B:179:THR:HG21	2.18	0.44
1:A:91:ARG:NH2	1:A:379:VAL:HG12	2.26	0.43
1:B:150:ASP:O	1:B:153:THR:OG1	2.28	0.42
1:A:102:VAL:HA	1:A:105:TYR:HD2	1.84	0.42
1:A:127:ASN:HD21	1:A:136:ALA:N	2.10	0.42
1:B:323:PRO:HB2	1:B:354:THR:HG22	2.01	0.42
1:B:339:TRP:CD1	2:D:9:ARG:HG3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:HG21	1:B:352:MET:HB3	2.01	0.42
1:B:292:ASN:O	1:B:292:ASN:ND2	2.48	0.42
1:A:164:SER:HA	1:A:310:LEU:HD21	2.03	0.41
1:A:57:VAL:HB	1:A:132:HIS:CD2	2.56	0.41
1:B:167:LEU:HD21	1:B:307:VAL:HA	2.01	0.41
1:B:164:SER:HA	1:B:310:LEU:HD21	2.02	0.41
1:A:127:ASN:ND2	1:A:136:ALA:H	2.12	0.40
1:A:219:HIS:HA	1:A:220:PRO:HD3	1.93	0.40
1:A:75:PHE:HB2	1:A:121:MET:SD	2.61	0.40
1:A:324:TYR:CZ	1:A:328:ARG:HD2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	298/335 (89%)	285 (96%)	13 (4%)	0	100	100
1	B	310/335 (92%)	296 (96%)	14 (4%)	0	100	100
2	C	4/10 (40%)	4 (100%)	0	0	100	100
2	D	4/10 (40%)	4 (100%)	0	0	100	100
All	All	616/690 (89%)	589 (96%)	27 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	264/290 (91%)	260 (98%)	4 (2%)	70	90
1	B	270/290 (93%)	268 (99%)	2 (1%)	87	95
2	C	6/7 (86%)	6 (100%)	0	100	100
2	D	6/7 (86%)	6 (100%)	0	100	100
All	All	546/594 (92%)	540 (99%)	6 (1%)	78	92

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

Mol	Chain	Res	Type
1	A	237	VAL
1	A	301	GLN
1	A	383	THR
1	A	384	LEU
1	B	133	HIS
1	B	292	ASN

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

Mol	Chain	Res	Type
1	A	127	ASN
1	A	211	GLN
1	A	219	HIS
1	A	239	GLN
1	A	301	GLN
1	A	365	ASN
1	A	370	ASN
1	B	127	ASN
1	B	211	GLN
1	B	219	HIS
1	B	239	GLN
1	B	270	GLN
1	B	365	ASN
1	B	370	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 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	GLY	A	1385	-	1,4,4	0.44	0	0,4,4	0.00	-
3	GLY	A	1386	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	A	1387	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	A	1388	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	A	1389	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	A	1390	-	1,4,4	0.44	0	0,4,4	0.00	-
3	GLY	A	1391	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	A	1392	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	B	1387	-	1,4,4	0.44	0	0,4,4	0.00	-
3	GLY	B	1388	-	1,4,4	0.43	0	0,4,4	0.00	-
3	GLY	B	1389	-	1,4,4	0.43	0	0,4,4	0.00	-

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	GLY	A	1385	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1386	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1387	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1388	-	-	0/0/2/2	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLY	A	1389	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1390	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1391	-	-	0/0/2/2	0/0/0/0
3	GLY	A	1392	-	-	0/0/2/2	0/0/0/0
3	GLY	B	1387	-	-	0/0/2/2	0/0/0/0
3	GLY	B	1388	-	-	0/0/2/2	0/0/0/0
3	GLY	B	1389	-	-	0/0/2/2	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1388	GLY	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	304/335 (90%)	0.78	44 (14%) 3 2	25, 63, 151, 300	0
1	B	314/335 (93%)	0.83	52 (16%) 2 1	32, 78, 163, 298	0
2	C	6/10 (60%)	1.30	1 (16%) 2 1	45, 52, 77, 84	0
2	D	6/10 (60%)	1.03	1 (16%) 2 1	57, 63, 87, 90	0
All	All	630/690 (91%)	0.81	98 (15%) 2 2	25, 69, 158, 300	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	99	GLN	10.3
1	B	372	VAL	7.5
1	B	98	LEU	7.0
1	A	91	ARG	6.3
1	A	269	HIS	5.6
1	B	376	PHE	5.6
1	A	376	PHE	5.5
1	B	371	LEU	5.4
1	B	200	LEU	4.6
1	B	197	SER	4.5
1	B	58	ASN	4.5
1	A	380	PHE	4.5
2	C	8	ARG	4.3
1	B	296	GLU	4.2
1	B	378	GLN	4.0
1	A	87	PHE	3.9
1	A	92	LYS	3.8
1	B	153	THR	3.6
1	A	244	MET	3.6
1	B	196	ALA	3.5
1	B	384	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	196	ALA	3.4
1	B	337	GLU	3.4
1	A	265	THR	3.3
1	A	171	ILE	3.2
1	A	248	PHE	3.2
1	B	244	MET	3.2
1	A	299	ARG	3.2
1	A	337	GLU	3.1
1	B	381	LEU	3.1
2	D	9	ARG	3.1
1	B	380	PHE	3.0
1	B	152	CYS	3.0
1	B	87	PHE	3.0
1	A	245	SER	3.0
1	A	360	VAL	3.0
1	B	198	ALA	3.0
1	A	224	VAL	2.9
1	B	199	LEU	2.9
1	B	100	SER	2.9
1	B	154	TYR	2.9
1	B	194	TRP	2.9
1	B	176	LYS	2.9
1	B	245	SER	2.9
1	B	55	LEU	2.8
1	A	137	PHE	2.8
1	A	210	LEU	2.8
1	B	155	ALA	2.8
1	B	374	ALA	2.8
1	B	370	ASN	2.8
1	B	373	SER	2.8
1	A	209	GLY	2.8
1	A	384	LEU	2.7
1	A	378	GLN	2.7
1	A	145	TYR	2.7
1	B	151	ALA	2.7
1	A	175	PHE	2.6
1	A	178	LYS	2.6
1	A	128	PHE	2.6
1	B	377	ARG	2.6
1	A	223	LEU	2.5
1	B	335	SER	2.5
1	A	383	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	339	TRP	2.5
1	A	199	LEU	2.5
1	B	102	VAL	2.5
1	A	242	THR	2.5
1	A	89	LEU	2.5
1	A	266	VAL	2.5
1	B	112	SER	2.5
1	B	50	GLY	2.5
1	B	295	ILE	2.4
1	B	119	LEU	2.4
1	A	334	ILE	2.4
1	B	203	PRO	2.3
1	B	201	ALA	2.3
1	A	197	SER	2.3
1	B	248	PHE	2.3
1	A	333	TYR	2.3
1	A	198	ALA	2.3
1	B	243	PHE	2.3
1	B	204	MET	2.3
1	A	157	ALA	2.2
1	B	228	ILE	2.2
1	A	247	LEU	2.2
1	A	135	TRP	2.2
1	A	136	ALA	2.2
1	A	246	PHE	2.2
1	B	52	ASN	2.1
1	B	195	LEU	2.1
1	A	225	CYS	2.1
1	A	243	PHE	2.1
1	B	193	ILE	2.1
1	A	200	LEU	2.1
1	B	158	LEU	2.1
1	A	183	ARG	2.1
1	B	49	SER	2.0
1	B	385	ALA	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	GLY	A	1387	5/5	0.69	0.67	7.53	92,92,92,125	0
3	GLY	A	1385	5/5	0.81	0.35	4.65	67,67,67,80	0
3	GLY	B	1387	5/5	0.78	0.53	3.55	79,94,94,94	0
3	GLY	A	1392	5/5	0.75	0.30	3.19	68,77,77,77	0
3	GLY	A	1389	5/5	0.87	0.31	0.81	89,89,89,107	0
3	GLY	B	1389	5/5	0.78	0.33	0.76	108,108,108,143	0
3	GLY	A	1386	5/5	0.72	0.20	-0.11	90,102,102,102	0
3	GLY	B	1388	5/5	0.83	0.42	-0.11	96,104,104,104	0
3	GLY	A	1388	5/5	0.89	0.23	-0.18	76,98,98,98	0
3	GLY	A	1391	5/5	0.66	0.45	-	89,89,89,101	0
3	GLY	A	1390	5/5	0.73	0.86	-	72,72,72,144	0

6.5 Other polymers [i](#)

There are no such residues in this entry.