



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:42 PM GMT

PDB ID : 5DMG
Title : X-RAY STRUCTURE OF THE FAB FRAGMENT OF THE ANTI TAU ANTIBODY RB86 IN COMPLEX WITH THE PHOSPHORYLATED TAU PEPTIDE (416-430)
Authors : Benz, J.; Lorenz, S.; Georges, G.; Jochner, A.; Goepfert, U.; Grueninger, F.; Bujotzek, A.
Deposited on : 2015-09-08
Resolution : 2.50 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
<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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20026982

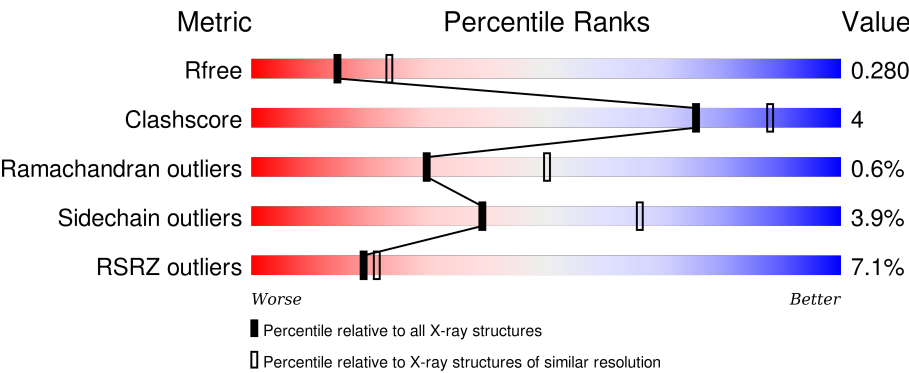
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	C	211	<div><div></div><div>87%8%</div></div>
1	E	211	<div><div>3%</div><div>80%12%6%</div></div>
1	H	211	<div><div>2%</div><div>89%7%</div></div>
2	D	219	<div><div></div><div>89%9%</div></div>
2	F	219	<div><div>32%</div><div>82%16%</div></div>

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Mol	Chain	Length	Quality of chain
2	L	219	<div><div>%</div><div><div></div><div>89%</div><div>9%</div><div>..</div></div></div>
3	P	15	<div><div>13%</div><div><div></div><div>60%</div><div>20%</div><div>20%</div></div></div>
3	X	15	<div><div>7%</div><div><div></div><div>40%</div><div>7%</div><div>53%</div></div></div>
3	Z	15	<div><div>7%</div><div><div></div><div>53%</div><div>7%</div><div>40%</div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9833 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 RB86 antibody Fab fragment heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	202	Total	C	N	O	S	0	0	0
			1490	948	244	293	5			
1	C	202	Total	C	N	O	S	0	0	0
			1490	948	244	293	5			
1	E	198	Total	C	N	O	S	0	0	0
			1464	933	239	287	5			

- Molecule 2 is a protein called RB86 antibody Fab fragment light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	216	Total	C	N	O	S	0	0	0
			1631	1019	272	333	7			
2	D	217	Total	C	N	O	S	0	0	0
			1640	1024	274	335	7			
2	F	215	Total	C	N	O	S	0	0	0
			1626	1015	272	332	7			

- Molecule 3 is a protein called Microtubule-associated protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Z	9	Total	C	N	O	P	0	0	0
			69	41	10	17	1			
3	P	12	Total	C	N	O	P	0	0	0
			90	53	13	22	1			
3	X	7	Total	C	N	O	P	0	0	0
			57	33	8	14	1			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	39	Total	O	0	0
			39	39		

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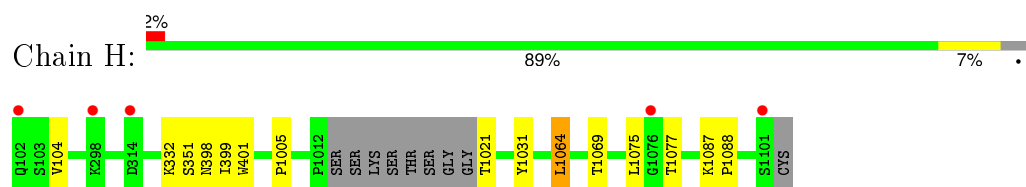
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	45	Total 45	O 45	0	0
4	C	54	Total 54	O 54	0	0
4	D	61	Total 61	O 61	0	0
4	E	51	Total 51	O 51	0	0
4	F	22	Total 22	O 22	0	0
4	Z	1	Total 1	O 1	0	0
4	P	1	Total 1	O 1	0	0
4	X	2	Total 2	O 2	0	0

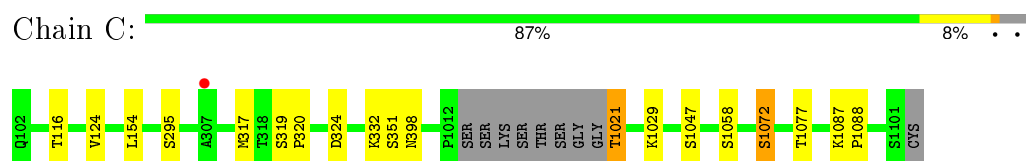
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

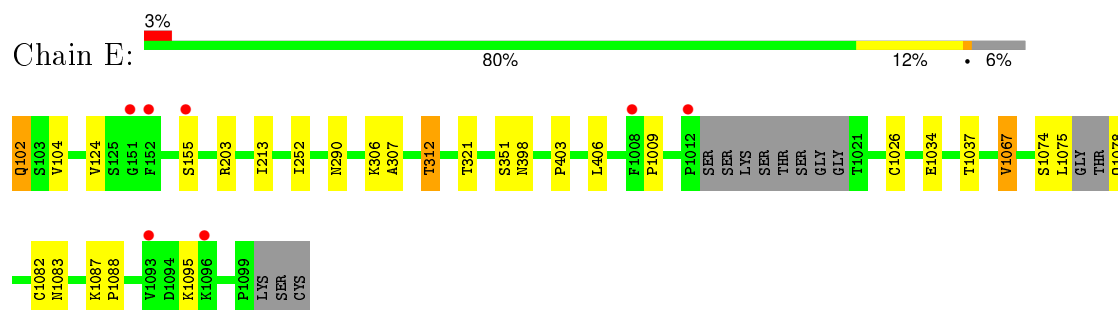
- Molecule 1: RB86 antibody Fab fragment heavy chain



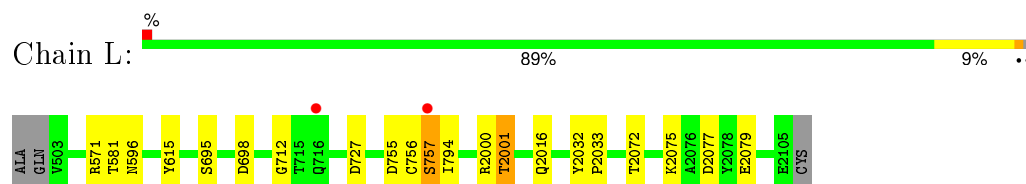
- Molecule 1: RB86 antibody Fab fragment heavy chain



- Molecule 1: RB86 antibody Fab fragment heavy chain

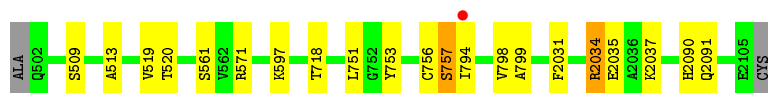


- Molecule 2: RB86 antibody Fab fragment light chain

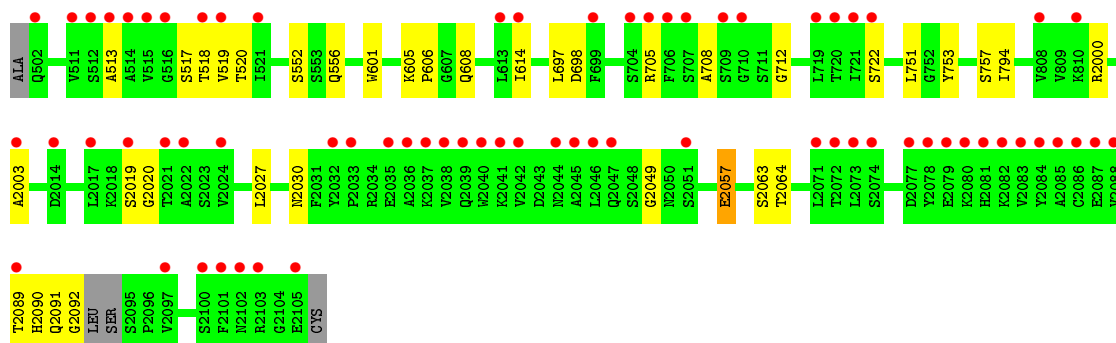
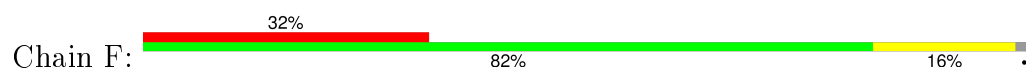


- Molecule 2: RB86 antibody Fab fragment light chain

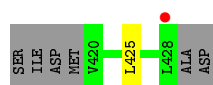




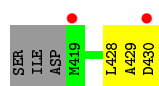
- Molecule 2: RB86 antibody Fab fragment light chain



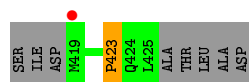
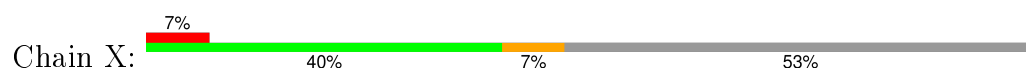
- Molecule 3: Microtubule-associated protein



- Molecule 3: Microtubule-associated protein



- Molecule 3: Microtubule-associated protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.51Å 163.05Å 70.93Å 90.00° 110.27° 90.00°	Depositor
Resolution (Å)	46.61 – 2.50 46.56 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.61-2.50) 99.9 (46.56-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.210 , 0.263 0.236 , 0.280	Depositor DCC
R_{free} test set	2509 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 49616 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9833	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SEP

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	C	0.33	0/1527	0.52	0/2089
1	E	0.37	0/1500	0.55	0/2052
1	H	0.32	0/1527	0.53	0/2089
2	D	0.36	0/1673	0.60	0/2276
2	F	0.32	0/1658	0.53	0/2254
2	L	0.35	0/1664	0.55	0/2264
3	P	0.33	0/79	0.52	0/105
3	X	0.39	0/46	0.52	0/59
3	Z	0.38	0/58	0.63	0/77
All	All	0.34	0/9732	0.55	0/13265

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

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	C	1490	0	1481	8	0
1	E	1464	0	1452	14	0
1	H	1490	0	1481	9	0
2	D	1640	0	1588	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1626	0	1571	16	0
2	L	1631	0	1580	17	0
3	P	90	0	82	5	0
3	X	57	0	50	1	0
3	Z	69	0	64	0	0
4	C	54	0	0	0	0
4	D	61	0	0	0	0
4	E	51	0	0	0	0
4	F	22	0	0	0	0
4	H	39	0	0	0	0
4	L	45	0	0	1	0
4	P	1	0	0	0	0
4	X	2	0	0	0	0
4	Z	1	0	0	0	0
All	All	9833	0	9349	73	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 4.

All (73) 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)
2:L:581:THR:HG22	2:L:596:ASN:N	1.55	1.15
2:L:581:THR:CG2	2:L:596:ASN:H	1.57	1.08
2:L:581:THR:HG22	2:L:596:ASN:H	0.71	0.78
1:E:1074:SER:HA	1:E:1075:LEU:C	2.07	0.74
2:F:753:TYR:OH	3:X:423:PRO:O	2.09	0.67
1:C:351:SER:OG	1:C:398:ASN:N	2.30	0.65
1:H:351:SER:OG	1:H:398:ASN:N	2.30	0.64
2:L:581:THR:CG2	2:L:596:ASN:N	2.30	0.61
2:D:2090:HIS:ND1	2:D:2091:GLN:O	2.34	0.59
2:F:2019:SER:N	2:F:2020:GLY:HA2	2.18	0.57
1:H:1021:THR:HG22	1:H:1069:THR:HG22	1.87	0.56
1:E:307:ALA:N	1:E:312:THR:O	2.33	0.55
2:D:756:CYS:O	2:D:757:SER:CB	2.54	0.55
2:L:581:THR:HG22	2:L:596:ASN:OD1	2.06	0.55
2:D:513:ALA:HB2	2:D:519:VAL:CG1	2.37	0.55
2:F:601:TRP:CD1	2:F:614:ILE:HD11	2.41	0.54
1:E:1067:VAL:HG21	2:F:2027:LEU:HD22	1.89	0.54
2:L:615:TYR:CZ	3:P:428:LEU:HB2	2.43	0.54
2:D:751:LEU:HD11	2:D:798:VAL:CG1	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:520:THR:HG23	2:D:718:THR:HG23	1.90	0.54
2:F:552:SER:OG	2:F:556:GLN:O	2.24	0.52
2:L:2000:ARG:HG2	2:L:2001:THR:N	2.25	0.52
2:F:513:ALA:HB2	2:F:519:VAL:HG11	1.92	0.51
1:H:332:LYS:HZ2	3:P:429:ALA:HB1	1.77	0.50
2:F:606:PRO:HG2	2:F:2057:GLU:HG3	1.93	0.50
1:E:307:ALA:HB3	1:E:312:THR:HB	1.92	0.50
1:E:351:SER:OG	1:E:398:ASN:N	2.44	0.49
2:F:614:ILE:HG22	2:F:697:LEU:HA	1.94	0.49
2:L:571:ARG:O	2:L:581:THR:C	2.50	0.49
2:L:2000:ARG:NH1	2:L:2001:THR:HG23	2.28	0.48
2:L:2075:LYS:O	2:L:2079:GLU:HG2	2.14	0.48
1:C:1021:THR:N	1:C:1072:SER:HG	2.12	0.47
2:L:756:CYS:O	2:L:757:SER:CB	2.61	0.47
3:P:429:ALA:O	3:P:430:ASP:C	2.53	0.47
2:D:597:LYS:HG2	2:D:753:TYR:CD1	2.50	0.47
2:F:705:ARG:HB3	2:F:722:SER:O	2.16	0.46
2:F:614:ILE:CD1	2:F:708:ALA:HB3	2.46	0.46
2:L:756:CYS:O	2:L:757:SER:HB3	2.15	0.46
2:L:755:ASP:HB3	2:L:794:ILE:HD12	1.97	0.46
2:L:615:TYR:CE1	3:P:428:LEU:HB2	2.51	0.46
2:L:2016:GLN:NE2	4:L:2201:HOH:O	2.47	0.46
1:E:203:ARG:HB3	1:E:213:ILE:HD11	1.98	0.45
2:L:596:ASN:O	2:L:596:ASN:CG	2.55	0.45
2:F:757:SER:HA	2:F:794:ILE:HA	1.72	0.44
2:F:2000:ARG:HG3	2:F:2063:SER:HB2	1.99	0.44
2:F:2030:ASN:OD1	2:F:2064:THR:HG21	2.18	0.44
1:H:104:VAL:HG22	1:H:399:ILE:HG22	1.99	0.44
1:E:102:GLN:HG2	1:E:102:GLN:O	2.18	0.43
1:C:124:VAL:HG21	1:C:154:LEU:HD21	2.00	0.43
2:F:517:SER:OG	2:F:518:THR:N	2.52	0.43
1:C:1087:LYS:N	1:C:1088:PRO:CD	2.81	0.43
1:E:1087:LYS:N	1:E:1088:PRO:CD	2.82	0.43
2:D:751:LEU:HD12	2:D:799:ALA:O	2.19	0.43
1:E:1009:PRO:HD3	1:E:1095:LYS:HE2	2.00	0.43
1:H:1064:LEU:HD12	1:H:1064:LEU:C	2.39	0.43
1:H:351:SER:O	1:H:401:TRP:NE1	2.52	0.42
1:C:332:LYS:O	1:C:351:SER:C	2.57	0.42
1:E:104:VAL:HG12	1:E:124:VAL:HA	2.02	0.42
1:C:317:MET:HE1	1:C:324:ASP:HB3	2.02	0.42
2:F:2000:ARG:NH1	2:F:2003:ALA:HB2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:319:SER:N	1:C:320:PRO:HD3	2.35	0.42
1:H:1005:PRO:HB3	1:H:1031:TYR:HB3	2.03	0.41
1:E:1074:SER:OG	1:E:1078:GLN:HB3	2.21	0.41
1:E:406:LEU:HD21	1:E:1034:GLU:HB2	2.02	0.41
2:D:2031:PHE:CE1	2:D:2034:ARG:HA	2.55	0.41
1:E:312:THR:HG22	1:E:312:THR:O	2.20	0.41
1:H:332:LYS:NZ	3:P:429:ALA:HB1	2.36	0.41
2:F:2090:HIS:O	2:F:2092:GLY:N	2.54	0.41
2:L:2032:TYR:CG	2:L:2033:PRO:HA	2.56	0.41
1:C:116:THR:O	1:C:320:PRO:HG2	2.21	0.41
1:H:1087:LYS:N	1:H:1088:PRO:CD	2.84	0.41
2:D:751:LEU:HD11	2:D:798:VAL:HG13	2.03	0.40
1:E:252:ILE:HG12	1:E:306:LYS:HD3	2.02	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	C	198/211 (94%)	197 (100%)	1 (0%)	0	100	100
1	E	192/211 (91%)	186 (97%)	6 (3%)	0	100	100
1	H	198/211 (94%)	191 (96%)	7 (4%)	0	100	100
2	D	215/219 (98%)	203 (94%)	11 (5%)	1 (0%)	34	55
2	F	211/219 (96%)	192 (91%)	16 (8%)	3 (1%)	14	24
2	L	214/219 (98%)	203 (95%)	9 (4%)	2 (1%)	21	37
3	P	9/15 (60%)	9 (100%)	0	0	100	100
3	X	4/15 (27%)	2 (50%)	1 (25%)	1 (25%)	0	0
3	Z	6/15 (40%)	6 (100%)	0	0	100	100
All	All	1247/1335 (93%)	1189 (95%)	51 (4%)	7 (1%)	30	50

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	2049	GLY
2	F	2091	GLN
2	L	757	SER
2	D	757	SER
2	F	712	GLY
2	L	712	GLY
3	X	423	PRO

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	C	170/177 (96%)	163 (96%)	7 (4%)	37	63
1	E	167/177 (94%)	156 (93%)	11 (7%)	21	38
1	H	170/177 (96%)	167 (98%)	3 (2%)	66	88
2	D	188/189 (100%)	181 (96%)	7 (4%)	41	68
2	F	186/189 (98%)	179 (96%)	7 (4%)	40	67
2	L	187/189 (99%)	181 (97%)	6 (3%)	46	74
3	P	9/12 (75%)	9 (100%)	0	100	100
3	X	6/12 (50%)	6 (100%)	0	100	100
3	Z	7/12 (58%)	6 (86%)	1 (14%)	4	7
All	All	1090/1134 (96%)	1048 (96%)	42 (4%)	39	66

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1064	LEU
1	H	1075	LEU
1	H	1077	THR
2	L	695	SER
2	L	698	ASP
2	L	727	ASP

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Mol	Chain	Res	Type
2	L	2001	THR
2	L	2072	THR
2	L	2077	ASP
1	C	295	SER
1	C	1021	THR
1	C	1029	LYS
1	C	1047	SER
1	C	1058	SER
1	C	1072	SER
1	C	1077	THR
2	D	509	SER
2	D	561	SER
2	D	571	ARG
2	D	794	ILE
2	D	2034	ARG
2	D	2035	GLU
2	D	2037	LYS
1	E	102	GLN
1	E	155	SER
1	E	290	ASN
1	E	312	THR
1	E	321	THR
1	E	403	PRO
1	E	1026	CYS
1	E	1037	THR
1	E	1067	VAL
1	E	1082	CYS
1	E	1083	ASN
2	F	520	THR
2	F	605	LYS
2	F	608	GLN
2	F	698	ASP
2	F	751	LEU
2	F	2057	GLU
2	F	2089	THR
3	Z	425	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	102	GLN
1	H	1090	ASN

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Mol	Chain	Res	Type
2	L	608	GLN
2	L	2039	GLN
2	L	2047	GLN
2	L	2052	GLN
1	C	290	ASN
2	D	502	GLN
2	D	551	GLN
2	D	596	ASN
2	D	608	GLN
2	D	716	GLN
2	D	2016	GLN
2	D	2030	ASN
2	D	2047	GLN
2	D	2091	GLN
1	E	290	ASN
2	F	2090	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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	SEP	P	422	3	7,9,10	0.71	0	8,12,14	1.67	3 (37%)
3	SEP	X	422	3	7,9,10	0.70	0	8,12,14	1.75	2 (25%)
3	SEP	Z	422	3	7,9,10	0.74	0	8,12,14	1.32	1 (12%)

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	SEP	P	422	3	-	0/5/8/10	0/0/0/0
3	SEP	X	422	3	-	0/5/8/10	0/0/0/0
3	SEP	Z	422	3	-	0/5/8/10	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Z	422	SEP	O-C-CA	-2.25	119.69	125.72
3	X	422	SEP	O-C-CA	-2.22	119.78	125.72
3	P	422	SEP	O-C-CA	-2.13	120.00	125.72
3	P	422	SEP	O3P-P-O2P	2.15	115.33	107.44
3	P	422	SEP	OG-CB-CA	2.80	110.70	108.26
3	X	422	SEP	OG-CB-CA	2.93	110.81	108.26

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	C	202/211 (95%)	0.09	1 (0%) 91 92	27, 36, 50, 66	0
1	E	198/211 (93%)	0.35	7 (3%) 48 53	21, 36, 53, 67	0
1	H	202/211 (95%)	0.12	5 (2%) 61 65	27, 37, 51, 62	0
2	D	217/219 (99%)	0.20	1 (0%) 91 92	22, 34, 51, 65	0
2	F	215/219 (98%)	1.58	70 (32%) 1 0	25, 47, 75, 84	0
2	L	216/219 (98%)	0.20	2 (0%) 85 88	23, 35, 52, 59	0
3	P	11/15 (73%)	0.84	2 (18%) 2 2	37, 42, 73, 76	0
3	X	6/15 (40%)	1.45	1 (16%) 2 2	72, 82, 100, 125	0
3	Z	8/15 (53%)	1.07	1 (12%) 5 5	58, 63, 73, 78	0
All	All	1275/1335 (95%)	0.44	90 (7%) 19 21	21, 38, 62, 125	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	2040	TRP	9.2
2	F	2019	SER	8.5
2	F	2045	ALA	8.2
2	F	2078	TYR	6.9
2	F	2083	VAL	6.8
3	X	419	MET	6.3
2	F	2084	TYR	6.2
2	F	706	PHE	5.3
2	F	518	THR	5.2
2	F	2044	ASN	5.0
2	F	2036	ALA	4.9
2	F	2101	PHE	4.8
2	F	721	ILE	4.3
2	F	2071	LEU	4.3
3	P	419	MET	4.2

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Mol	Chain	Res	Type	RSRZ
2	F	2046	LEU	4.0
2	F	2082	LYS	3.9
2	F	2022	ALA	3.9
2	F	2073	LEU	3.9
2	F	2085	ALA	3.8
2	F	2105	GLU	3.6
2	F	2081	HIS	3.6
2	F	513	ALA	3.4
2	F	2024	VAL	3.4
2	F	2102	ASN	3.4
2	F	2086	CYS	3.4
2	F	2077	ASP	3.4
2	F	2103	ARG	3.4
3	P	430	ASP	3.3
2	F	514	ALA	3.3
2	F	2079	GLU	3.2
2	F	2037	LYS	3.2
2	F	2072	THR	3.2
2	F	719	LEU	3.2
2	F	516	GLY	3.1
1	E	152	PHE	3.1
2	F	512	SER	3.0
1	E	155	SER	3.0
2	F	519	VAL	2.9
2	F	2088	VAL	2.9
2	F	2087	GLU	2.9
2	F	2033	PRO	2.9
2	F	2038	VAL	2.9
2	F	722	SER	2.9
1	E	1012	PRO	2.9
2	F	2021	THR	2.9
2	F	2074	SER	2.8
2	F	2039	GLN	2.8
2	F	613	LEU	2.8
2	F	2041	LYS	2.8
2	F	521	ILE	2.7
2	F	2017	LEU	2.7
1	E	1096	LYS	2.7
2	F	705	ARG	2.7
1	H	1101	SER	2.7
2	F	704	SER	2.6
2	F	2042	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
2	F	2014	ASP	2.6
2	F	709	SER	2.6
2	L	757	SER	2.6
2	F	707	SER	2.6
1	E	1093	VAL	2.6
2	F	808	VAL	2.5
1	C	307	ALA	2.5
2	F	511	VAL	2.5
1	H	298	LYS	2.5
2	F	2097	VAL	2.5
2	F	2035	GLU	2.5
1	E	1008	PHE	2.4
2	F	720	THR	2.4
2	F	2032	TYR	2.3
2	F	515	VAL	2.3
1	H	1076	GLY	2.2
1	H	102	GLN	2.2
1	E	151	GLY	2.2
1	H	314	ASP	2.2
2	F	614	ILE	2.2
2	F	2047	GLN	2.2
2	F	699	PHE	2.2
2	F	710	GLY	2.2
2	L	716	GLN	2.1
3	Z	428	LEU	2.1
2	F	2003	ALA	2.1
2	F	2089	THR	2.1
2	F	810	LYS	2.1
2	F	2100	SER	2.1
2	F	2080	LYS	2.1
2	F	2051	SER	2.1
2	F	502	GLN	2.0
2	D	794	ILE	2.0

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

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	SEP	P	422	10/11	0.95	0.11	-	39,42,46,49	0
3	SEP	Z	422	10/11	0.95	0.17	-	58,60,66,66	0
3	SEP	X	422	10/11	0.91	0.18	-	77,78,81,83	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.