



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 04:00 PM EDT

PDB ID : 6QEP / pdb_00006qep
Title : EngBF DARPin Fusion 4b H14
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Deposited on : 2019-01-08
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

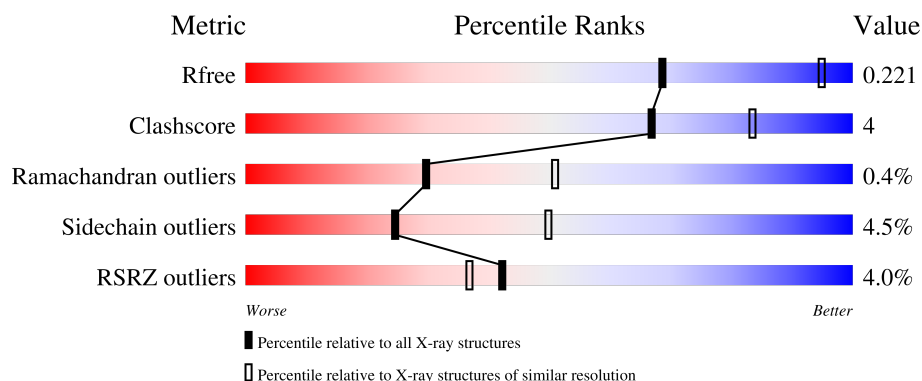
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	1354	 4% 84% 13% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10896 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.

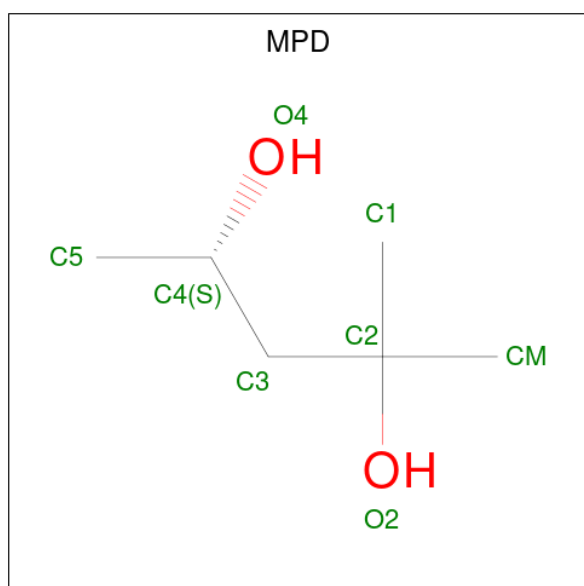
- Molecule 1 is a protein called PEGA domain-containing protein,EngBF-DARPin Fusion 4b H14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1341	10299	6412	1770	2091	26	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	334	GLY	-	expression tag	UNP A0A374RF80
A	335	PRO	-	expression tag	UNP A0A374RF80
A	336	LEU	-	expression tag	UNP A0A374RF80
A	337	GLY	-	expression tag	UNP A0A374RF80
A	338	SER	-	expression tag	UNP A0A374RF80
A	339	MET	-	expression tag	UNP A0A374RF80
A	1309	ARG	GLN	conflict	UNP A0A374RF80

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	Mn	0	0
			4	4		

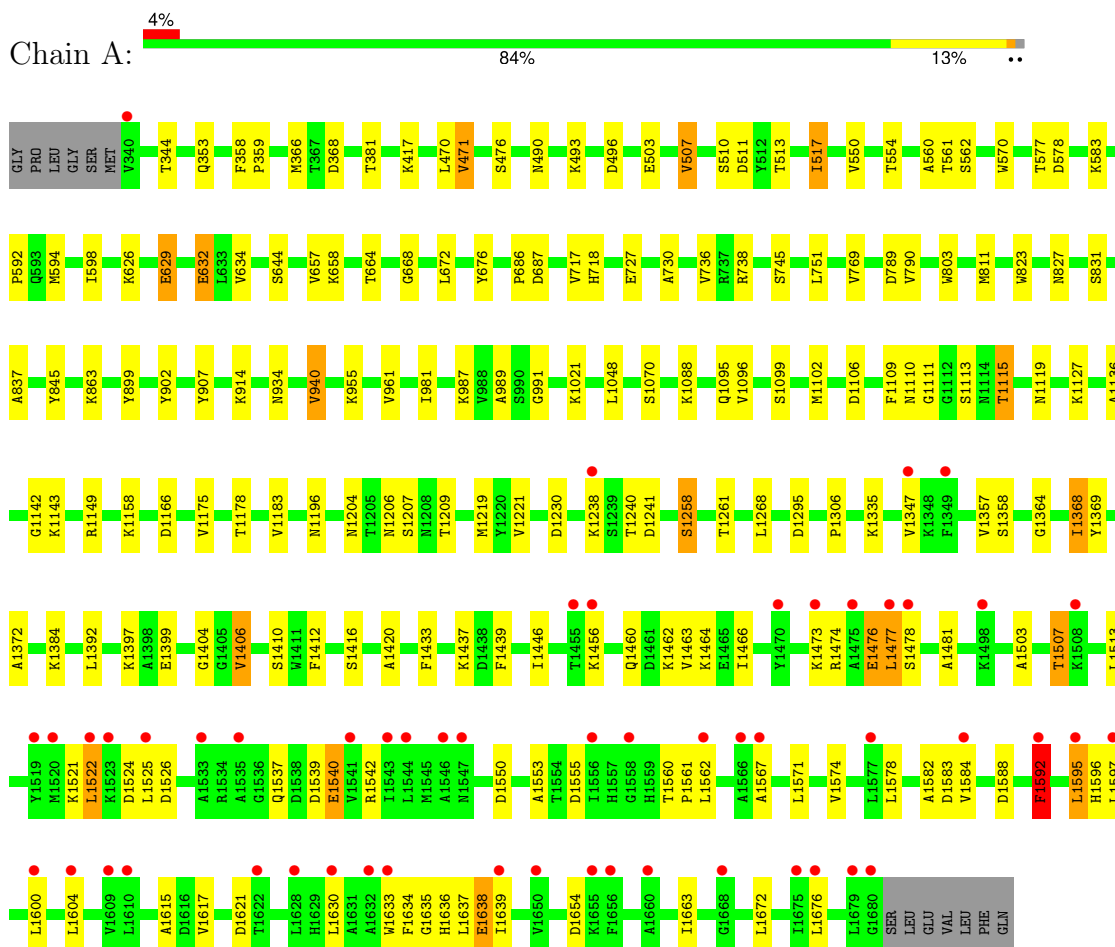
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	573	Total	O	0	0
			573	573		

3 Residue-property plots

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: PEGA domain-containing protein,EngBF-DARPin Fusion 4b H14



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	192.69Å 192.69Å 123.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.28 – 2.60 46.28 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.28-2.60) 99.9 (46.28-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.61Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.177 , 0.215 0.180 , 0.221	Depositor DCC
R_{free} test set	4017 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 79.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.024 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10896	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MES, MPD, MN

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	0.86	0/10498	1.33	48/14239 (0.3%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1592	PHE	CA-CB-CG	9.43	123.23	113.80
1	A	1588	ASP	CA-C-N	7.57	131.31	120.79
1	A	1588	ASP	C-N-CA	7.57	131.31	120.79
1	A	934	ASN	CA-CB-CG	7.26	119.86	112.60
1	A	1241	ASP	CA-CB-CG	6.87	119.47	112.60
1	A	629	GLU	CA-C-N	6.83	124.54	120.24
1	A	629	GLU	C-N-CA	6.83	124.54	120.24
1	A	1464	LYS	CA-C-N	6.73	129.19	120.44
1	A	1464	LYS	C-N-CA	6.73	129.19	120.44
1	A	727	GLU	CB-CA-C	-6.58	99.49	110.68
1	A	789	ASP	CA-C-N	6.45	133.58	121.97
1	A	789	ASP	C-N-CA	6.45	133.58	121.97
1	A	578	ASP	CA-CB-CG	6.40	119.00	112.60
1	A	1621	ASP	CA-CB-CG	6.33	118.93	112.60
1	A	1460	GLN	CA-C-N	6.08	128.71	120.38
1	A	1460	GLN	C-N-CA	6.08	128.71	120.38
1	A	1221	VAL	N-CA-CB	5.88	121.10	111.45
1	A	513	THR	N-CA-C	-5.83	105.00	111.36
1	A	1295	ASP	N-CA-C	-5.78	98.46	108.52
1	A	1654	ASP	CA-CB-CG	5.67	118.27	112.60
1	A	1555	ASP	CA-CB-CG	5.59	118.19	112.60
1	A	1306	PRO	CA-C-N	5.56	128.29	120.28
1	A	1306	PRO	C-N-CA	5.56	128.29	120.28
1	A	1456	LYS	N-CA-C	-5.50	105.28	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	837	ALA	N-CA-C	-5.49	105.63	112.93
1	A	368	ASP	CA-CB-CG	5.46	118.06	112.60
1	A	1110	ASN	CA-CB-CG	5.45	118.05	112.60
1	A	1113	SER	CA-C-N	5.44	128.12	120.28
1	A	1113	SER	C-N-CA	5.44	128.12	120.28
1	A	751	LEU	N-CA-C	-5.41	104.94	112.45
1	A	1521	LYS	CA-C-N	5.32	129.68	120.58
1	A	1521	LYS	C-N-CA	5.32	129.68	120.58
1	A	1588	ASP	CA-CB-CG	5.30	117.90	112.60
1	A	1404	GLY	CA-C-N	5.28	124.99	119.92
1	A	1404	GLY	C-N-CA	5.28	124.99	119.92
1	A	687	ASP	CA-CB-CG	5.24	117.84	112.60
1	A	1578	LEU	CA-C-N	5.23	127.29	120.28
1	A	1578	LEU	C-N-CA	5.23	127.29	120.28
1	A	1433	PHE	CA-C-N	5.21	125.86	120.03
1	A	1433	PHE	C-N-CA	5.21	125.86	120.03
1	A	1521	LYS	N-CA-C	-5.16	103.51	110.68
1	A	1206	ASN	CA-CB-CG	5.13	117.73	112.60
1	A	991	GLY	CA-C-N	5.09	129.74	123.12
1	A	991	GLY	C-N-CA	5.09	129.74	123.12
1	A	803	TRP	CA-C-N	5.01	127.00	120.28
1	A	803	TRP	C-N-CA	5.01	127.00	120.28
1	A	823	TRP	CA-C-N	5.01	125.64	120.03
1	A	823	TRP	C-N-CA	5.01	125.64	120.03

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	10299	0	9868	74	0
2	A	8	0	14	0	0
3	A	12	0	13	0	0
4	A	4	0	0	0	0
5	A	573	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10896	0	9895	74	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 (74) 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:471:VAL:HG13	1:A:517:ILE:HD11	1.55	0.89
1:A:1149:ARG:HH12	1:A:1230:ASP:HA	1.49	0.78
1:A:1142:GLY:HA2	1:A:1240:THR:HA	1.71	0.71
1:A:1571:LEU:HA	1:A:1574:VAL:HG22	1.74	0.68
1:A:1115:THR:HG22	1:A:1119:ASN:OD1	1.92	0.68
1:A:1462:LYS:O	1:A:1466:ILE:HG12	1.93	0.68
1:A:1364:GLY:HA2	1:A:1392:LEU:HD13	1.76	0.65
1:A:507:VAL:HG22	1:A:510:SER:HB2	1.78	0.65
1:A:1178:THR:HG22	1:A:1183:VAL:HA	1.80	0.63
1:A:1143:LYS:HG2	1:A:1238:LYS:O	1.99	0.62
1:A:1550:ASP:HB3	1:A:1553:ALA:HB2	1.81	0.62
1:A:1175:VAL:HG12	1:A:1219:MET:HE3	1.82	0.60
1:A:1595:LEU:HD11	1:A:1615:ALA:HB1	1.83	0.59
1:A:1106:ASP:OD2	1:A:1115:THR:HG21	2.01	0.59
1:A:570:TRP:CD2	1:A:592:PRO:HB3	2.39	0.57
1:A:507:VAL:CG2	1:A:510:SER:HB2	2.35	0.57
1:A:1567:ALA:HB2	1:A:1597:LEU:HB3	1.88	0.54
1:A:955:LYS:HG2	1:A:961:VAL:HG22	1.89	0.54
1:A:1600:LEU:HG	1:A:1630:LEU:HB3	1.90	0.53
1:A:1473:LYS:HB3	1:A:1476:GLU:HG2	1.91	0.53
1:A:1335:LYS:HB2	1:A:1416:SER:HB3	1.92	0.52
1:A:1503:ALA:O	1:A:1507:THR:HG22	2.08	0.51
1:A:1463:VAL:HG22	1:A:1513:LEU:HD11	1.94	0.50
1:A:1478:SER:HB3	1:A:1481:ALA:HB3	1.93	0.50
1:A:344:THR:HG22	1:A:353:GLN:HG3	1.95	0.49
1:A:1111:GLY:HA3	1:A:1115:THR:HG23	1.94	0.49
1:A:1109:PHE:O	1:A:1136:ALA:HB3	2.12	0.49
1:A:554:THR:HA	1:A:562:SER:O	2.13	0.49
1:A:1166:ASP:OD2	1:A:1209:THR:HG21	2.12	0.49
1:A:476:SER:HB3	1:A:560:ALA:HB2	1.95	0.48
1:A:1368:ILE:HD11	1:A:1420:ALA:C	2.38	0.48
1:A:1503:ALA:O	1:A:1507:THR:CG2	2.61	0.48
1:A:1637:LEU:HD12	1:A:1672:LEU:HA	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1537:GLN:HB3	1:A:1540:GLU:HB2	1.96	0.47
1:A:1143:LYS:HG3	1:A:1240:THR:HB	1.96	0.47
1:A:1149:ARG:HH12	1:A:1230:ASP:CA	2.23	0.47
1:A:1347:VAL:HG11	1:A:1446:ILE:HD13	1.95	0.47
1:A:632:GLU:O	1:A:668:GLY:HA3	2.15	0.47
1:A:626:LYS:O	1:A:629:GLU:HG2	2.14	0.47
1:A:1477:LEU:HD22	1:A:1522:LEU:HD13	1.97	0.46
1:A:1562:LEU:HD11	1:A:1582:ALA:HB1	1.97	0.46
1:A:658:LYS:HD3	1:A:1096:VAL:HG13	1.97	0.46
1:A:827:ASN:HB3	1:A:831:SER:HB2	1.97	0.46
1:A:634:VAL:HA	1:A:914:LYS:HG3	1.97	0.45
1:A:1539:ASP:HA	1:A:1542:ARG:HE	1.82	0.45
1:A:1604:LEU:HD22	1:A:1638:GLU:HB2	1.97	0.45
1:A:736:VAL:HG12	1:A:738:ARG:HG3	1.99	0.45
1:A:1592:PHE:HB2	1:A:1596:HIS:CG	2.52	0.44
1:A:664:THR:HA	1:A:907:TYR:OH	2.18	0.44
1:A:718:HIS:CD2	1:A:718:HIS:C	2.96	0.44
1:A:899:TYR:O	1:A:902:TYR:HB3	2.18	0.44
1:A:470:LEU:HD22	1:A:594:MET:HE2	2.00	0.43
1:A:1149:ARG:NH1	1:A:1230:ASP:HA	2.26	0.43
1:A:769:VAL:HG23	1:A:811:MET:HG2	2.00	0.43
1:A:1357:VAL:O	1:A:1399:GLU:HA	2.17	0.43
1:A:981:ILE:HD12	1:A:989:ALA:HB3	2.01	0.43
1:A:1048:LEU:HD11	1:A:1088:LYS:HG2	2.01	0.43
1:A:1560:THR:HB	1:A:1561:PRO:HD2	2.00	0.43
1:A:358:PHE:CG	1:A:359:PRO:HD2	2.54	0.43
1:A:1369:TYR:CD2	1:A:1439:PHE:HB2	2.53	0.43
1:A:831:SER:O	1:A:863:LYS:HE2	2.19	0.42
1:A:490:ASN:HA	1:A:845:TYR:O	2.21	0.41
1:A:577:THR:OG1	1:A:583:LYS:HG2	2.20	0.41
1:A:676:TYR:CD1	1:A:717:VAL:HB	2.55	0.41
1:A:366:MET:HE3	1:A:598:ILE:HG21	2.03	0.41
1:A:1099:SER:HB3	1:A:1102:MET:HB2	2.02	0.41
1:A:1372:ALA:O	1:A:1412:PHE:HA	2.21	0.41
1:A:1158:LYS:NZ	1:A:1258:SER:HB3	2.35	0.41
1:A:1635:GLY:HA2	1:A:1672:LEU:HD11	2.03	0.41
1:A:686:PRO:HD2	1:A:730:ALA:HB3	2.01	0.41
1:A:490:ASN:ND2	1:A:493:LYS:HD2	2.36	0.40
1:A:1600:LEU:HD11	1:A:1630:LEU:HD22	2.03	0.40
1:A:511:ASP:HA	1:A:550:VAL:O	2.22	0.40
1:A:1196:ASN:HB3	1:A:1204:ASN:HA	2.03	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	1339/1354 (99%)	1271 (95%)	62 (5%)	6 (0%)	30 51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	790	VAL
1	A	940	VAL
1	A	1524	ASP
1	A	1583	ASP
1	A	1638	GLU
1	A	1406	VAL

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	1093/1104 (99%)	1044 (96%)	49 (4%)	24 50

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	381	THR
1	A	417	LYS

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Mol	Chain	Res	Type
1	A	471	VAL
1	A	496	ASP
1	A	503	GLU
1	A	507	VAL
1	A	517	ILE
1	A	561	THR
1	A	632	GLU
1	A	644	SER
1	A	657	VAL
1	A	672	LEU
1	A	745	SER
1	A	940	VAL
1	A	987	LYS
1	A	1021	LYS
1	A	1070	SER
1	A	1095	GLN
1	A	1115	THR
1	A	1127	LYS
1	A	1207	SER
1	A	1258	SER
1	A	1261	THR
1	A	1268	LEU
1	A	1358	SER
1	A	1368	ILE
1	A	1384	LYS
1	A	1397	LYS
1	A	1406	VAL
1	A	1410	SER
1	A	1437	LYS
1	A	1474	ARG
1	A	1476	GLU
1	A	1477	LEU
1	A	1507	THR
1	A	1522	LEU
1	A	1525	LEU
1	A	1526	ASP
1	A	1540	GLU
1	A	1584	VAL
1	A	1592	PHE
1	A	1595	LEU
1	A	1617	VAL
1	A	1633	TRP

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Mol	Chain	Res	Type
1	A	1634	PHE
1	A	1636	HIS
1	A	1639	ILE
1	A	1663	ILE
1	A	1676	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	388	ASN
1	A	443	ASN
1	A	467	ASN
1	A	492	GLN
1	A	555	GLN
1	A	793	ASN
1	A	813	ASN
1	A	1025	GLN
1	A	1256	GLN
1	A	1452	GLN
1	A	1484	GLN
1	A	1570	HIS
1	A	1587	ASN
1	A	1613	ASN
1	A	1677	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
2	MPD	A	1701	-	7,7,7	0.60	0	9,10,10	0.31	0
3	MES	A	1702	-	12,12,12	0.68	0	15,16,16	0.40	0

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
2	MPD	A	1701	-	-	2/5/5/5	-
3	MES	A	1702	-	-	2/6/14/14	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1702	MES	C8-C7-N4-C5
3	A	1702	MES	C8-C7-N4-C3
2	A	1701	MPD	O2-C2-C3-C4
2	A	1701	MPD	C1-C2-C3-C4

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1341/1354 (99%)	-0.06	54 (4%) 42 37	43, 74, 213, 255	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1477	LEU	4.2
1	A	1610	LEU	4.0
1	A	1622	THR	3.7
1	A	1520	MET	3.7
1	A	1522	LEU	3.7
1	A	1577	LEU	3.6
1	A	1656	PHE	3.5
1	A	1650	VAL	3.4
1	A	1655	LYS	3.3
1	A	1478	SER	3.3
1	A	1592	PHE	3.3
1	A	1556	ILE	3.2
1	A	1595	LEU	3.1
1	A	1675	ILE	3.1
1	A	1628	LEU	3.1
1	A	1584	VAL	3.1
1	A	1633	TRP	3.0
1	A	1525	LEU	2.9
1	A	1597	LEU	2.9
1	A	1600	LEU	2.9
1	A	1347	VAL	2.8
1	A	1679	LEU	2.8
1	A	340	VAL	2.7
1	A	1535	ALA	2.7
1	A	1604	LEU	2.6
1	A	1630	LEU	2.6
1	A	1470	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	1609	VAL	2.5
1	A	1546	ALA	2.5
1	A	1455	THR	2.5
1	A	1349	PHE	2.5
1	A	1680	GLY	2.4
1	A	1519	TYR	2.4
1	A	1562	LEU	2.4
1	A	1632	ALA	2.4
1	A	1543	ILE	2.3
1	A	1541	VAL	2.3
1	A	1639	ILE	2.3
1	A	1523	LYS	2.3
1	A	1498	LYS	2.3
1	A	1473	LYS	2.2
1	A	1456	LYS	2.2
1	A	1238	LYS	2.2
1	A	1668	GLY	2.2
1	A	1660	ALA	2.2
1	A	1475	ALA	2.1
1	A	1547	ASN	2.1
1	A	1544	LEU	2.1
1	A	1566	ALA	2.1
1	A	1533	ALA	2.1
1	A	1567	ALA	2.1
1	A	1508	LYS	2.1
1	A	1676	LEU	2.0
1	A	1558	GLY	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 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
2	MPD	A	1701	8/8	0.91	0.21	101,103,105,106	0
3	MES	A	1702	12/12	0.94	0.13	88,100,108,109	0
4	MN	A	1706	1/1	0.98	0.04	83,83,83,83	0
4	MN	A	1703	1/1	0.99	0.04	91,91,91,91	0
4	MN	A	1705	1/1	1.00	0.02	78,78,78,78	0
4	MN	A	1704	1/1	1.00	0.01	72,72,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.