



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:41 PM GMT

PDB ID : 1BZH
Title : Cyclic peptide inhibitor of human PTP1B
Authors : Groves, M.R.; Yao, Z.J.; Burke Jr., T.R.; Barford, D.
Deposited on : 1998-10-28
Resolution : 2.10 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

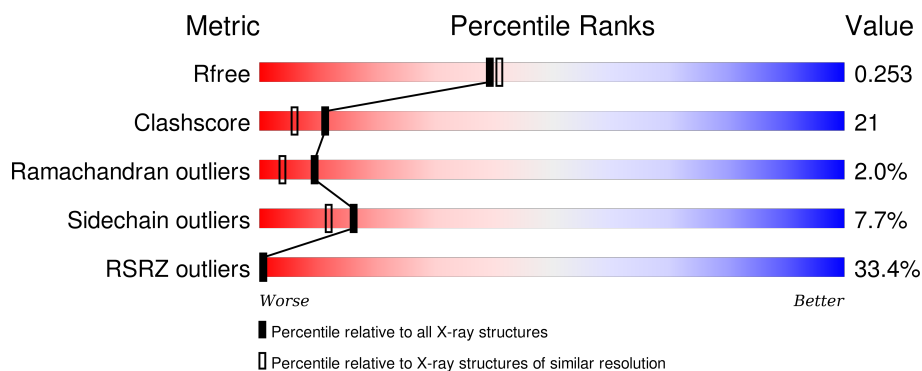
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	298	
2	I	7	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2583 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 (PROTEIN-TYROSINE-PHOSPHATASE 1B).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2427	1535	418	458	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	252	ASP	GLU	CONFLICT	UNP P18031
A	265	GLU	ASP	CONFLICT	UNP P18031

- Molecule 2 is a protein called PROTEIN (PROTEIN-TYROSINE-PHOSPHATASE 1B INHIBITOR).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	7	Total	C	F	N	O	S	0	0
			68	39	1	8	19	1		

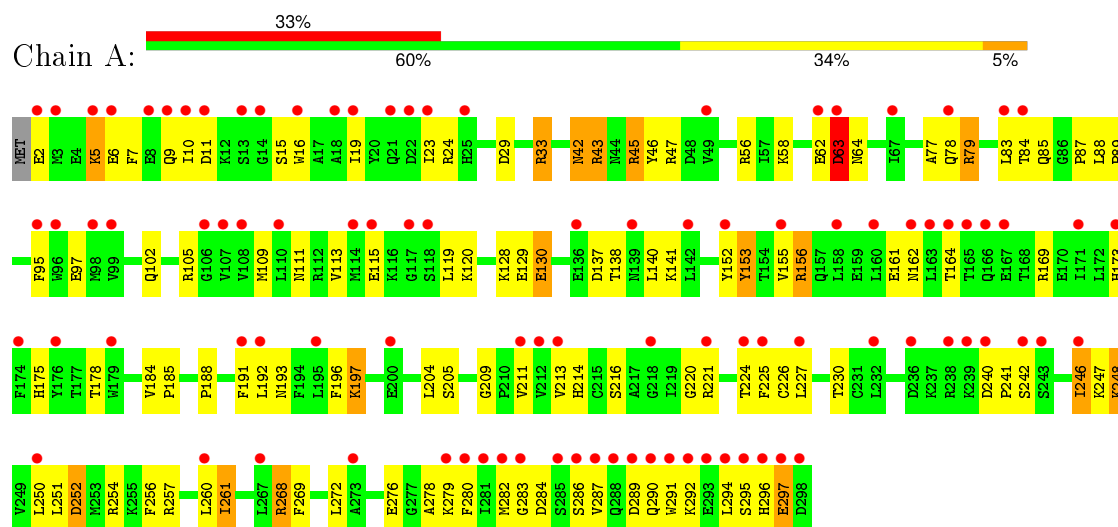
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	88	Total	O	0	0
			88	88		

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 ($\text{RSRZ} > 2$). 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 (PROTEIN-TYROSINE-PHOSPHATASE 1B)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.69Å 86.82Å 52.00Å 90.00° 96.91° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10 24.74 – 2.10	Depositor EDS
% Data completeness (in resolution range)	78.1 (8.00-2.10) 71.3 (24.74-2.10)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.56 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.196 , 0.262 0.204 , 0.253	Depositor DCC
R_{free} test set	1371 reflections (10.29%)	DCC
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 84.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 14569 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2583	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.67% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.54	0/2482	1.41	26/3344 (0.8%)
2	I	1.79	1/36 (2.8%)	2.41	2/46 (4.3%)
All	All	0.58	1/2518 (0.0%)	1.43	28/3390 (0.8%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	401	ASP	C-N	5.17	1.46	1.34

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	ARG	NE-CZ-NH1	9.30	124.95	120.30
1	A	169	ARG	NE-CZ-NH2	-9.10	115.75	120.30
1	A	43	ARG	NE-CZ-NH1	-8.96	115.82	120.30
1	A	257	ARG	NE-CZ-NH2	-8.65	115.97	120.30
1	A	156	ARG	NE-CZ-NH2	8.33	124.47	120.30
1	A	256	PHE	CB-CG-CD2	-8.22	115.05	120.80
1	A	79	ARG	NE-CZ-NH2	-7.91	116.35	120.30
1	A	252	ASP	CB-CG-OD2	7.70	125.23	118.30
1	A	63	ASP	CB-CG-OD1	6.83	124.45	118.30
1	A	137	ASP	CB-CG-OD2	-6.17	112.75	118.30
1	A	268	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	129	GLU	OE1-CD-OE2	-5.94	116.17	123.30
1	A	24	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	A	254	ARG	NE-CZ-NH2	5.77	123.19	120.30
1	A	153	TYR	CB-CG-CD2	5.76	124.46	121.00
1	A	29	ASP	CB-CG-OD1	5.75	123.48	118.30
1	A	33	ARG	NE-CZ-NH2	5.65	123.12	120.30
2	I	404	GLU	OE1-CD-OE2	-5.63	116.54	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ARG	NE-CZ-NH2	-5.63	117.49	120.30
1	A	33	ARG	CA-CB-CG	5.62	125.77	113.40
1	A	153	TYR	CB-CG-CD1	-5.46	117.73	121.00
1	A	169	ARG	NH1-CZ-NH2	5.45	125.39	119.40
1	A	105	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	79	ARG	CD-NE-CZ	5.09	130.72	123.60
1	A	268	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	A	252	ASP	N-CA-CB	5.04	119.68	110.60
1	A	45	ARG	NE-CZ-NH1	-5.03	117.78	120.30
2	I	401	ASP	C-N-CA	-5.03	109.13	121.70

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	A	2427	0	2381	93	0
2	I	68	0	43	18	0
3	A	88	0	0	3	0
All	All	2583	0	2424	102	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 21.

All (102) 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:97:GLU:HB2	1:A:138:THR:HG21	1.32	1.11
1:A:45:ARG:H	1:A:85:GLN:HE22	1.16	0.94
1:A:46:TYR:HA	2:I:404:GLU:HG3	1.56	0.86
1:A:220:GLY:HA3	2:I:405:FLT:O2	1.78	0.83
1:A:287:VAL:HG21	1:A:290:GLN:HB2	1.62	0.82
1:A:291:TRP:HA	1:A:294:LEU:HD13	1.61	0.81
1:A:83:LEU:HD21	1:A:226:CYS:SG	2.21	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:ARG:H	1:A:85:GLN:NE2	1.82	0.77
2:I:401:ASP:CG	2:I:402:ALA:H	1.90	0.74
1:A:155:VAL:HG21	1:A:197:LYS:NZ	2.04	0.71
2:I:401:ASP:CG	2:I:402:ALA:N	2.45	0.70
1:A:152:TYR:CE1	1:A:295:SER:HB3	2.28	0.69
1:A:97:GLU:HB2	1:A:138:THR:CG2	2.16	0.68
1:A:88:LEU:HB3	1:A:89:PRO:HD2	1.75	0.68
2:I:403:ASP:OD1	2:I:404:GLU:OE1	2.11	0.67
1:A:248:LYS:HE3	1:A:251:LEU:HB2	1.76	0.67
1:A:46:TYR:CD1	2:I:405:FLT:HB3	2.32	0.65
1:A:250:LEU:CD1	1:A:261:ILE:HD13	2.28	0.64
1:A:5:LYS:HD2	1:A:6:GLU:N	2.12	0.64
1:A:250:LEU:HD13	1:A:261:ILE:HD13	1.80	0.63
1:A:47:ARG:HG2	2:I:404:GLU:HG2	1.81	0.62
1:A:102:GLN:O	1:A:209:GLY:HA3	1.99	0.62
2:I:401:ASP:N	2:I:401:ASP:OD2	2.34	0.61
1:A:152:TYR:OH	1:A:193:ASN:HB3	2.00	0.61
1:A:193:ASN:HD22	1:A:291:TRP:HB2	1.66	0.60
1:A:155:VAL:HG21	1:A:197:LYS:HZ1	1.67	0.59
1:A:153:TYR:OH	1:A:197:LYS:HG3	2.03	0.58
1:A:77:ALA:CB	1:A:230:THR:HG23	2.34	0.58
1:A:56:ARG:NH2	1:A:58:LYS:HZ2	2.03	0.57
1:A:294:LEU:H	1:A:294:LEU:HD12	1.70	0.56
1:A:83:LEU:HD22	1:A:213:VAL:HB	1.87	0.56
1:A:46:TYR:CZ	2:I:405:FLT:HD1	2.41	0.55
1:A:109:MET:HG3	1:A:214:HIS:CE1	2.42	0.55
2:I:403:ASP:C	2:I:404:GLU:OE1	2.45	0.54
1:A:111:ASN:HD21	1:A:214:HIS:CE1	2.25	0.54
1:A:287:VAL:CG2	1:A:290:GLN:HB2	2.34	0.54
1:A:140:LEU:HD23	1:A:162:ASN:HA	1.90	0.53
1:A:56:ARG:NH2	1:A:58:LYS:NZ	2.56	0.53
1:A:155:VAL:HG21	1:A:197:LYS:HZ3	1.74	0.53
1:A:241:PRO:HG3	1:A:282:MET:SD	2.48	0.53
1:A:188:PRO:HD3	1:A:272:LEU:HD13	1.90	0.53
1:A:23:ILE:HD11	1:A:247:LYS:HG2	1.90	0.52
1:A:16:TRP:NE1	1:A:268:ARG:HD2	2.23	0.52
1:A:47:ARG:H	2:I:404:GLU:HG3	1.72	0.52
2:I:405:FLT:C	2:I:405:FLT:CD2	2.87	0.52
1:A:280:PHE:CE1	1:A:290:GLN:HB3	2.46	0.51
1:A:119:LEU:HD22	3:A:330:HOH:O	2.11	0.51
1:A:287:VAL:HG11	1:A:290:GLN:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:PHE:HZ	1:A:225:PHE:HA	1.77	0.49
1:A:120:LYS:HB3	1:A:216:SER:HB2	1.94	0.49
1:A:77:ALA:O	1:A:78:GLN:HB2	2.13	0.48
1:A:33:ARG:HA	3:A:347:HOH:O	2.13	0.48
1:A:6:GLU:HG2	1:A:10:ILE:CD1	2.44	0.48
1:A:77:ALA:HB2	1:A:230:THR:HG23	1.93	0.48
1:A:16:TRP:CE2	1:A:268:ARG:HD2	2.49	0.47
1:A:184:VAL:HB	1:A:185:PRO:CD	2.45	0.47
1:A:47:ARG:NH1	2:I:404:GLU:HG2	2.30	0.47
1:A:138:THR:HG22	1:A:138:THR:O	2.14	0.47
1:A:276:GLU:HG3	1:A:280:PHE:HE2	1.80	0.46
1:A:2:GLU:O	1:A:5:LYS:HE3	2.15	0.46
1:A:47:ARG:H	2:I:404:GLU:CG	2.29	0.46
1:A:42:ASN:ND2	1:A:43:ARG:HG3	2.32	0.45
2:I:403:ASP:HB3	2:I:404:GLU:CD	2.37	0.45
1:A:162:ASN:OD1	1:A:164:THR:HB	2.17	0.45
1:A:46:TYR:CG	2:I:405:FLT:HB3	2.52	0.45
1:A:287:VAL:HG11	1:A:290:GLN:HG3	1.98	0.45
1:A:191:PHE:CZ	1:A:225:PHE:HA	2.52	0.44
1:A:56:ARG:HH21	1:A:58:LYS:NZ	2.14	0.44
1:A:63:ASP:HB2	1:A:64:ASN:H	1.64	0.44
1:A:193:ASN:ND2	1:A:291:TRP:HE3	2.16	0.44
1:A:278:ALA:O	1:A:282:MET:HB2	2.17	0.44
1:A:115:GLU:HB2	1:A:120:LYS:HG3	2.00	0.44
1:A:204:LEU:HD21	1:A:211:VAL:HG11	2.00	0.44
1:A:111:ASN:O	1:A:175:HIS:NE2	2.45	0.44
1:A:7:PHE:CE1	1:A:268:ARG:HG3	2.52	0.44
1:A:193:ASN:ND2	1:A:291:TRP:HB2	2.30	0.44
1:A:227:LEU:HD21	1:A:246:ILE:HD11	1.99	0.44
1:A:87:PRO:HG3	1:A:95:PHE:CD2	2.52	0.44
1:A:56:ARG:HH21	1:A:58:LYS:HZ2	1.64	0.44
1:A:156:ARG:NH2	3:A:307:HOH:O	2.51	0.44
1:A:156:ARG:HB2	1:A:173:HIS:HB3	2.00	0.43
1:A:192:LEU:O	1:A:196:PHE:CD2	2.72	0.43
1:A:45:ARG:N	1:A:85:GLN:HE22	1.98	0.43
1:A:83:LEU:HD12	1:A:260:LEU:HD22	2.00	0.43
1:A:188:PRO:HA	1:A:269:PHE:CZ	2.54	0.42
1:A:240:ASP:OD2	1:A:242:SER:OG	2.33	0.42
1:A:88:LEU:HB3	1:A:89:PRO:CD	2.47	0.42
1:A:152:TYR:HA	1:A:178:THR:HG21	2.01	0.42
1:A:260:LEU:O	1:A:261:ILE:HB	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:VAL:CB	1:A:290:GLN:HB2	2.49	0.42
1:A:5:LYS:HD3	1:A:9:GLN:OE1	2.19	0.42
2:I:402:ALA:O	2:I:403:ASP:C	2.58	0.42
2:I:404:GLU:OE1	2:I:404:GLU:N	2.53	0.42
1:A:191:PHE:CZ	1:A:224:THR:HG22	2.55	0.42
1:A:287:VAL:HG11	1:A:290:GLN:CG	2.50	0.41
1:A:211:VAL:HG13	1:A:213:VAL:HG23	2.02	0.41
1:A:84:THR:O	1:A:214:HIS:HB2	2.20	0.41
1:A:6:GLU:HG2	1:A:10:ILE:HD12	2.03	0.41
1:A:280:PHE:CD1	1:A:290:GLN:HB3	2.55	0.41
1:A:2:GLU:HG3	1:A:5:LYS:HB3	2.03	0.41
1:A:128:LYS:HB3	1:A:130:GLU:OE2	2.21	0.40
1:A:141:LYS:HE3	1:A:161:GLU:OE1	2.22	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	295/298 (99%)	273 (92%)	16 (5%)	6 (2%)	9	4
2	I	3/7 (43%)	2 (67%)	1 (33%)	0	100	100
All	All	298/305 (98%)	275 (92%)	17 (6%)	6 (2%)	9	4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	ASP
1	A	296	HIS
1	A	297	GLU
1	A	261	ILE
1	A	283	GLY

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Mol	Chain	Res	Type
1	A	292	LYS

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	270/271 (100%)	250 (93%)	20 (7%)	17	13
2	I	4/4 (100%)	3 (75%)	1 (25%)	1	0
All	All	274/275 (100%)	253 (92%)	21 (8%)	16	12

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	11	ASP
1	A	15	SER
1	A	19	ILE
1	A	42	ASN
1	A	62	GLU
1	A	63	ASP
1	A	79	ARG
1	A	113	VAL
1	A	130	GLU
1	A	197	LYS
1	A	205	SER
1	A	246	ILE
1	A	248	LYS
1	A	252	ASP
1	A	279	LYS
1	A	284	ASP
1	A	286	SER
1	A	289	ASP
1	A	297	GLU
2	I	404	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	85	GLN
1	A	111	ASN
1	A	123	GLN
1	A	157	GLN
1	A	193	ASN
1	A	288	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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$
2	FLT	I	405	2	11,20,21	1.07	1 (9%)	12,28,30	0.77	0
2	AEA	I	407	2	9,9,10	2.31	2 (22%)	7,10,12	3.00	6 (85%)

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	FLT	I	405	2	-	0/6/23/25	0/1/1/1
2	AEA	I	407	2	-	0/7/9/10	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	405	FLT	CA-N	-2.44	1.40	1.47
2	I	407	AEA	C4-S1	2.74	1.83	1.81
2	I	407	AEA	C2-N2	6.03	1.44	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	I	407	AEA	C1-C3-S1	-4.47	102.27	112.84
2	I	407	AEA	C5-C4-S1	-3.23	105.83	112.71
2	I	407	AEA	O3-C2-N2	-3.08	118.52	123.08
2	I	407	AEA	C3-C1-C2	2.28	114.91	109.54
2	I	407	AEA	C2-C1-N1	2.51	119.91	108.84
2	I	407	AEA	C1-C2-N2	3.33	121.97	116.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	405	FLT	5	0

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	A	297/298 (99%)	1.48	97 (32%)  	45, 61, 100, 100	0
2	I	5/7 (71%)	4.15	4 (80%)  	84, 86, 89, 94	0
All	All	302/305 (99%)	1.52	101 (33%)  	45, 61, 100, 100	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	287	VAL	11.9
1	A	291	TRP	8.7
2	I	401	ASP	8.7
1	A	289	ASP	7.2
1	A	110	LEU	6.6
1	A	294	LEU	6.4
1	A	296	HIS	5.9
1	A	293	GLU	5.5
1	A	288	GLN	5.2
1	A	117	GLY	5.2
1	A	295	SER	5.1
1	A	9	GLN	5.0
1	A	95	PHE	4.9
2	I	403	ASP	4.8
1	A	292	LYS	4.8
1	A	83	LEU	4.6
1	A	96	TRP	4.5
1	A	164	THR	4.5
1	A	191	PHE	4.4
1	A	2	GLU	4.4
1	A	298	ASP	4.3
1	A	11	ASP	4.1
1	A	62	GLU	4.1
1	A	108	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	165	THR	4.0
1	A	285	SER	4.0
1	A	286	SER	3.9
1	A	107	VAL	3.9
1	A	290	GLN	3.8
1	A	67	ILE	3.8
1	A	179	TRP	3.8
1	A	84	THR	3.7
1	A	212	VAL	3.7
2	I	404	GLU	3.7
1	A	18	ALA	3.7
1	A	22	ASP	3.7
1	A	166	GLN	3.6
1	A	8	GLU	3.5
1	A	162	ASN	3.5
1	A	19	ILE	3.5
1	A	225	PHE	3.5
1	A	280	PHE	3.5
1	A	260	LEU	3.4
1	A	267	LEU	3.4
1	A	297	GLU	3.4
1	A	213	VAL	3.4
1	A	158	LEU	3.4
1	A	250	LEU	3.3
1	A	99	VAL	3.3
1	A	171	ILE	3.2
1	A	5	LYS	3.2
1	A	13	SER	3.2
1	A	10	ILE	3.2
1	A	118	SER	3.1
1	A	160	LEU	3.1
1	A	246	ILE	3.0
1	A	152	TYR	3.0
1	A	14	GLY	3.0
1	A	192	LEU	3.0
1	A	195	LEU	2.9
1	A	243	SER	2.9
1	A	142	LEU	2.9
1	A	239	LYS	2.9
1	A	242	SER	2.8
1	A	211	VAL	2.7
1	A	232	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	167	GLU	2.6
1	A	224	THR	2.6
1	A	115	GLU	2.6
1	A	106	GLY	2.6
1	A	3	MET	2.6
1	A	240	ASP	2.6
1	A	155	VAL	2.5
1	A	283	GLY	2.5
1	A	238	ARG	2.5
1	A	136	GLU	2.5
1	A	21	GLN	2.4
1	A	98	MET	2.4
1	A	139	ASN	2.4
1	A	279	LYS	2.4
1	A	25	HIS	2.4
1	A	227	LEU	2.4
1	A	114	MET	2.3
1	A	174	PHE	2.3
1	A	23	ILE	2.3
1	A	173	HIS	2.3
1	A	16	TRP	2.3
1	A	236	ASP	2.3
1	A	273	ALA	2.2
1	A	163	LEU	2.2
1	A	218	GLY	2.2
2	I	402	ALA	2.2
1	A	282	MET	2.2
1	A	49	VAL	2.2
1	A	6	GLU	2.2
1	A	221	ARG	2.2
1	A	200	GLU	2.2
1	A	281	ILE	2.1
1	A	176	TYR	2.1
1	A	63	ASP	2.1
1	A	78	GLN	2.1

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
2	AEA	I	407	10/11	0.79	0.40	-	88,90,93,93	0
2	FLT	I	405	20/21	0.81	0.19	-	72,76,80,80	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.