



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:19 PM UTC

PDB ID : 6D9W / pdb_00006d9w
Title : Crystal structure of Deinococcus radiodurans MntH, an Nramp-family transition metal transporter, in the inward-open apo state
Authors : Gaudet, R.; Bane, L.B.; Weihofen, W.A.; Singharoy, A.; Zimanyi, C.M.; Bozzi, A.T.
Deposited on : 2018-04-30
Resolution : 3.94 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
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

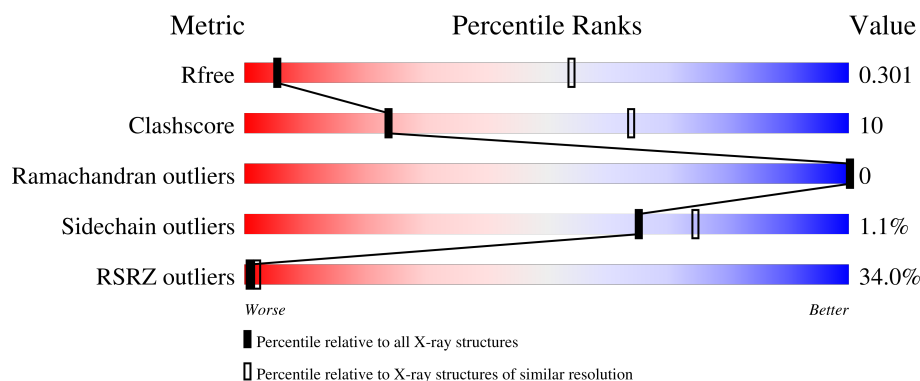
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.94 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	1022 (4.12-3.76)
Clashscore	190562	1000 (4.10-3.78)
Ramachandran outliers	187476	1009 (4.12-3.76)
Sidechain outliers	187428	1002 (4.12-3.76)
RSRZ outliers	180081	1022 (4.12-3.76)

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	420	43% 70%15%14%
2	H	213	15% 70%28%..
3	L	213	24% 76%24%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5764 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 Divalent metal cation transporter MntH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2545	1673	410	446	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	MET	-	initiating methionine	UNP Q9RTP8
A	18	HIS	-	expression tag	UNP Q9RTP8
A	19	HIS	-	expression tag	UNP Q9RTP8
A	20	HIS	-	expression tag	UNP Q9RTP8
A	21	HIS	-	expression tag	UNP Q9RTP8
A	22	HIS	-	expression tag	UNP Q9RTP8
A	23	HIS	-	expression tag	UNP Q9RTP8
A	24	HIS	-	expression tag	UNP Q9RTP8
A	25	HIS	-	expression tag	UNP Q9RTP8
A	168	HIS	GLN	engineered mutation	UNP Q9RTP8
A	169	HIS	LYS	engineered mutation	UNP Q9RTP8
A	251	TYR	GLU	engineered mutation	UNP Q9RTP8
A	252	TYR	GLU	engineered mutation	UNP Q9RTP8
A	253	TYR	LYS	engineered mutation	UNP Q9RTP8
A	398	HIS	ARG	engineered mutation	UNP Q9RTP8
A	399	HIS	ARG	engineered mutation	UNP Q9RTP8

- Molecule 2 is a protein called Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	211	Total	C	N	O	S	0	0	0
			1568	992	251	317	8			

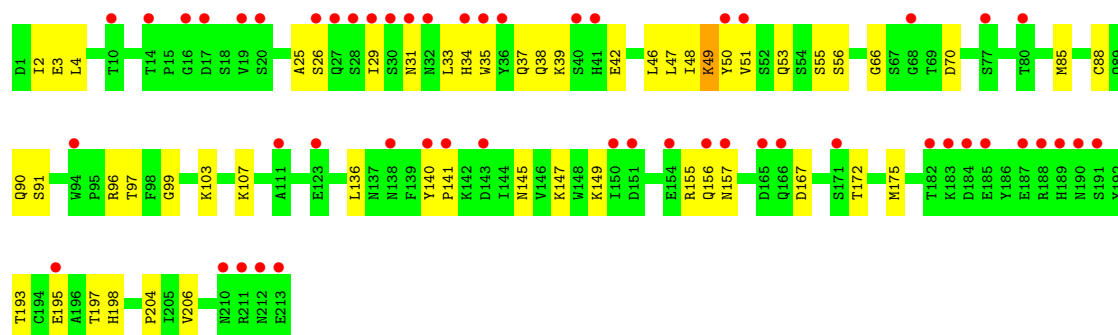
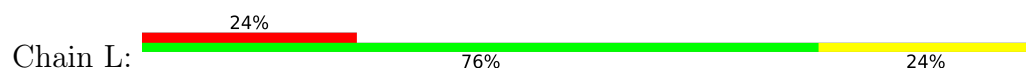
- Molecule 3 is a protein called Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total 1648	C 1019	N 282	O 341	S 6	0	0	0

- Molecule 4 is OSMIUM ION (CCD ID: OS) (formula: Os).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	Os 2	0	0
4	H	1	Total 1	Os 1	0	0

● Molecule 3: Fab Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	113.13Å 132.08Å 221.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.47 – 3.94 46.47 – 3.94	Depositor EDS
% Data completeness (in resolution range)	76.0 (46.47-3.94) 75.1 (46.47-3.94)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.43 (at 4.00Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.271 , 0.309 0.266 , 0.301	Depositor DCC
R_{free} test set	1179 reflections (7.86%)	wwPDB-VP
Wilson B-factor (Å ²)	-4.7	Xtriage
Anisotropy	5.658	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.74	EDS
Total number of atoms	5764	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% 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:
OS

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	0.13	0/2595	0.31	0/3560
2	H	0.14	0/1610	0.39	0/2207
3	L	0.11	0/1685	0.34	0/2286
All	All	0.13	0/5890	0.34	0/8053

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	1
3	L	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	98	GLU	Peptide
3	L	49	LYS	Peptide

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	2545	0	2539	40	0
2	H	1568	0	1532	43	0
3	L	1648	0	1565	35	0
4	A	2	0	0	0	0
4	H	1	0	0	0	0
All	All	5764	0	5636	115	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 10.

All (115) 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:393:LEU:HD21	1:A:412:THR:HG23	1.58	0.82
2:H:49:VAL:HG12	2:H:59:TYR:HA	1.70	0.74
1:A:84:MET:HE1	1:A:422:ILE:HD11	1.77	0.66
2:H:99:PRO:HB2	2:H:100:PRO:HD3	1.76	0.65
1:A:50:ALA:HB1	1:A:271:ALA:HB2	1.77	0.65
1:A:215:PRO:HB3	1:A:432:GLU:HG2	1.80	0.63
2:H:40:PRO:HB2	2:H:43:LYS:HB2	1.81	0.62
2:H:20:ILE:HD11	2:H:110:LEU:HD22	1.81	0.62
3:L:3:GLU:HA	3:L:97:THR:HG21	1.81	0.61
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.82	0.61
1:A:84:MET:HE2	1:A:229:VAL:HG21	1.82	0.61
1:A:288:GLY:O	2:H:101:TYR:OH	2.13	0.60
1:A:76:TRP:CE2	1:A:210:PRO:HB3	2.37	0.60
3:L:85:MET:HE2	3:L:103:LYS:HD2	1.82	0.60
2:H:20:ILE:HG21	2:H:108:THR:HG21	1.84	0.60
2:H:153:VAL:HG22	2:H:198:VAL:HG22	1.84	0.59
2:H:53:SER:HA	2:H:71:LYS:HE3	1.84	0.59
2:H:124:PRO:HB2	2:H:211:LEU:HD13	1.86	0.58
1:A:103:LEU:HD23	1:A:235:TYR:HE1	1.69	0.57
1:A:85:ALA:HB2	1:A:229:VAL:HG22	1.86	0.57
3:L:51:VAL:HG11	3:L:66:GLY:HA3	1.87	0.56
3:L:149:LYS:HB2	3:L:193:THR:HB	1.87	0.56
1:A:80:ALA:O	1:A:84:MET:HG2	2.06	0.55
3:L:91:SER:HA	3:L:96:ARG:HD2	1.88	0.55
3:L:33:LEU:O	3:L:34:HIS:ND1	2.39	0.55
2:H:24:VAL:HG21	2:H:29:LEU:HB2	1.89	0.55
1:A:226:GLY:HA3	1:A:382:CYS:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:4:LEU:HD23	3:L:25:ALA:HA	1.90	0.54
2:H:48:LEU:HG	2:H:49:VAL:HG22	1.87	0.54
3:L:38:GLN:HB3	3:L:85:MET:HB2	1.89	0.54
2:H:93:TYR:HE2	2:H:110:LEU:HD23	1.73	0.54
1:A:75:LEU:HD23	1:A:279:LEU:HD23	1.90	0.54
2:H:6:GLU:H	2:H:106:GLN:HE22	1.54	0.54
2:H:154:THR:OG1	2:H:197:SER:OG	2.25	0.54
2:H:51:ILE:HD13	2:H:57:THR:HG22	1.91	0.53
1:A:55:MET:HG3	1:A:183:VAL:HG21	1.89	0.53
3:L:88:CYS:O	3:L:99:GLY:N	2.39	0.53
2:H:12:VAL:HG21	2:H:18:LEU:HD23	1.90	0.53
1:A:81:ALA:HB1	1:A:228:THR:OG1	2.09	0.52
3:L:35:TRP:HB2	3:L:48:ILE:HB	1.91	0.52
1:A:62:THR:HG22	1:A:375:ILE:HG13	1.91	0.52
2:H:157:SER:HB3	2:H:194:VAL:HG13	1.92	0.51
2:H:47:TRP:CG	3:L:96:ARG:HB2	2.45	0.51
2:H:164:VAL:HG22	2:H:182:VAL:HG23	1.93	0.51
1:A:179:VAL:HG11	1:A:326:SER:HB2	1.93	0.50
1:A:416:TRP:HA	1:A:416:TRP:CE3	2.46	0.50
2:H:184:VAL:HG11	2:H:194:VAL:HG21	1.92	0.50
2:H:98:GLU:O	2:H:101:TYR:N	2.45	0.50
3:L:2:ILE:HG23	3:L:26:SER:HB2	1.92	0.50
1:A:160:VAL:HG11	1:A:325:LEU:HD13	1.94	0.50
3:L:2:ILE:HG21	3:L:29:ILE:HD11	1.94	0.49
3:L:4:LEU:HD11	3:L:90:GLN:HB3	1.94	0.49
1:A:382:CYS:O	1:A:385:VAL:HG12	2.13	0.49
1:A:179:VAL:HG13	1:A:322:ALA:HB1	1.95	0.49
2:H:12:VAL:HG23	2:H:112:VAL:HG22	1.95	0.49
2:H:90:THR:HG22	2:H:112:VAL:H	1.78	0.49
1:A:193:GLN:NE2	1:A:306:PRO:HD2	2.27	0.48
2:H:2:VAL:HA	2:H:26:GLY:HA3	1.95	0.48
3:L:145:ASN:HB3	3:L:197:THR:OG1	2.13	0.48
3:L:195:GLU:HG2	3:L:206:VAL:HG22	1.96	0.47
2:H:24:VAL:HG11	2:H:29:LEU:HD13	1.96	0.47
3:L:39:LYS:HB2	3:L:42:GLU:HG3	1.97	0.47
2:H:70:SER:HB2	2:H:79:PHE:HB2	1.96	0.47
3:L:197:THR:HG22	3:L:204:PRO:HB3	1.96	0.47
1:A:65:GLU:CD	1:A:295:GLY:HA3	2.40	0.47
3:L:31:ASN:O	3:L:50:TYR:HB3	2.15	0.47
1:A:363:VAL:HG11	1:A:373:VAL:HG11	1.97	0.46
2:H:12:VAL:O	2:H:112:VAL:HA	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:60:ASN:OD1	2:H:61:SER:N	2.49	0.46
1:A:65:GLU:OE2	1:A:295:GLY:HA3	2.16	0.46
2:H:35:HIS:NE2	2:H:101:TYR:OH	2.49	0.46
3:L:51:VAL:HG11	3:L:66:GLY:CA	2.46	0.46
1:A:132:LEU:HD13	1:A:360:ALA:HA	1.98	0.45
2:H:139:LEU:HD23	2:H:211:LEU:HG	1.98	0.45
1:A:76:TRP:CD2	1:A:210:PRO:HB3	2.52	0.45
2:H:99:PRO:CB	2:H:100:PRO:HD3	2.47	0.45
2:H:144:LYS:NZ	2:H:172:GLN:OE1	2.48	0.45
1:A:88:ILE:HD13	1:A:389:LEU:HD21	1.98	0.44
1:A:402:MET:HB2	1:A:406:VAL:HG22	1.98	0.44
1:A:297:LEU:HA	1:A:300:ALA:HB3	1.98	0.44
1:A:62:THR:HB	1:A:224:ILE:HG12	1.99	0.44
3:L:49:LYS:HG2	3:L:53:GLN:O	2.16	0.44
1:A:227:ALA:HB2	1:A:378:GLN:HB3	1.99	0.44
3:L:55:SER:OG	3:L:56:SER:N	2.50	0.44
1:A:357:MET:HE3	1:A:357:MET:HB3	1.89	0.43
3:L:147:LYS:NZ	3:L:156:GLN:HG2	2.33	0.43
3:L:157:ASN:OD1	3:L:157:ASN:N	2.51	0.43
1:A:222:VAL:HG13	1:A:382:CYS:SG	2.59	0.43
1:A:230:MET:HG3	1:A:233:VAL:H	1.83	0.43
3:L:107:LYS:HA	3:L:140:TYR:OH	2.19	0.43
2:H:162:SER:HA	2:H:183:THR:O	2.19	0.43
3:L:141:PRO:HD2	3:L:198:HIS:HE1	1.84	0.43
1:A:128:MET:SD	1:A:359:PRO:HG2	2.59	0.43
3:L:155:ARG:NH2	3:L:157:ASN:O	2.50	0.42
2:H:52:TRP:HB3	2:H:56:SER:HB2	2.00	0.42
2:H:24:VAL:HB	2:H:27:PHE:CE1	2.55	0.42
1:A:222:VAL:HA	1:A:225:ILE:HG22	2.02	0.42
3:L:70:ASP:OD1	3:L:70:ASP:N	2.53	0.42
2:H:63:LEU:HB3	2:H:67:LEU:HD13	2.02	0.42
2:H:100:PRO:O	3:L:46:LEU:HD22	2.20	0.42
3:L:46:LEU:HD23	3:L:55:SER:OG	2.20	0.42
1:A:193:GLN:OE1	1:A:281:VAL:HG11	2.20	0.41
1:A:281:VAL:O	1:A:285:THR:OG1	2.30	0.41
2:H:178:MET:HE2	2:H:178:MET:HB3	1.86	0.41
3:L:167:ASP:HB2	3:L:172:THR:O	2.21	0.41
2:H:157:SER:HB2	2:H:194:VAL:HG22	2.03	0.41
3:L:141:PRO:HD2	3:L:198:HIS:CE1	2.56	0.41
3:L:136:LEU:HD12	3:L:175:MET:SD	2.61	0.41
1:A:54:TYR:OH	1:A:82:ASN:ND2	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:19:SER:HA	2:H:80:LEU:O	2.21	0.41
3:L:147:LYS:HA	3:L:147:LYS:HD2	1.88	0.40
1:A:379:VAL:O	1:A:383:PHE:HD1	2.04	0.40
2:H:210:LYS:HG2	2:H:212:GLU:HG3	2.02	0.40
2:H:99:PRO:HB2	2:H:100:PRO:CD	2.50	0.40
2:H:35:HIS:CE1	2:H:50:VAL:HB	2.57	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	354/420 (84%)	345 (98%)	9 (2%)	0	100	100
2	H	207/213 (97%)	195 (94%)	12 (6%)	0	100	100
3	L	211/213 (99%)	201 (95%)	10 (5%)	0	100	100
All	All	772/846 (91%)	741 (96%)	31 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	241/326 (74%)	237 (98%)	4 (2%)	53	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	185/187 (99%)	182 (98%)	3 (2%)	55	70
3	L	192/193 (100%)	192 (100%)	0	100	100
All	All	618/706 (88%)	611 (99%)	7 (1%)	65	74

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	303	THR
1	A	411	PHE
1	A	416	TRP
2	H	50	VAL
2	H	141	CYS
2	H	196	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	426	ASN
3	L	41	HIS
3	L	53	GLN
3	L	161	ASN

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 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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.

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	362/420 (86%)	2.34	182 (50%) 0 1	55, 121, 168, 221	0
2	H	211/213 (99%)	0.94	33 (15%) 5 8	13, 59, 88, 126	0
3	L	213/213 (100%)	1.38	52 (24%) 2 4	40, 72, 114, 140	0
All	All	786/846 (92%)	1.70	267 (33%) 1 2	13, 84, 156, 221	0

All (267) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	109	GLU	13.7
1	A	108	ARG	11.7
1	A	299	THR	10.1
1	A	232	HIS	9.2
3	L	16	GLY	8.1
1	A	165	LEU	7.4
1	A	200	ASP	6.9
1	A	233	VAL	6.8
3	L	17	ASP	6.7
1	A	235	TYR	6.6
1	A	236	LEU	6.6
1	A	110	ARG	6.6
1	A	105	GLU	6.2
1	A	401	VAL	6.2
1	A	409	ARG	6.1
1	A	388	ALA	6.0
1	A	107	ILE	5.9
1	A	308	LEU	5.8
3	L	211	ARG	5.8
1	A	295	GLY	5.7
1	A	292	GLU	5.7
1	A	410	SER	5.5
2	H	75	LYS	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	261	ARG	5.3
1	A	408	ARG	5.3
1	A	402	MET	5.2
3	L	188	ARG	5.1
1	A	398	HIS	5.1
1	A	407	THR	5.1
1	A	231	PRO	5.0
1	A	112	PRO	4.9
1	A	204	VAL	4.9
1	A	54	TYR	4.8
1	A	303	THR	4.8
3	L	210	ASN	4.7
2	H	40	PRO	4.7
1	A	296	ASP	4.6
1	A	136	LEU	4.6
3	L	157	ASN	4.6
1	A	103	LEU	4.5
1	A	384	GLY	4.5
1	A	436	GLY	4.5
1	A	189	ALA	4.5
2	H	213	PRO	4.4
1	A	53	ALA	4.3
1	A	102	ASN	4.3
1	A	293	ASN	4.3
1	A	365	LEU	4.3
3	L	77	SER	4.2
1	A	195	VAL	4.2
1	A	257	VAL	4.2
1	A	198	ARG	4.2
1	A	152	TRP	4.2
1	A	302	GLN	4.2
3	L	212	ASN	4.1
1	A	404	ALA	4.1
1	A	215	PRO	4.1
1	A	196	LEU	4.0
1	A	358	LEU	4.0
1	A	397	ALA	4.0
1	A	162	PHE	4.0
3	L	26	SER	4.0
1	A	352	ARG	4.0
1	A	415	GLY	3.9
1	A	140	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
3	L	27	GLN	3.9
3	L	34	HIS	3.9
1	A	307	LEU	3.8
3	L	68	GLY	3.8
3	L	154	GLU	3.8
3	L	184	ASP	3.8
1	A	201	LEU	3.7
1	A	114	PRO	3.7
3	L	190	ASN	3.7
2	H	8	GLY	3.7
1	A	403	GLY	3.7
1	A	405	LEU	3.7
1	A	399	HIS	3.7
1	A	151	PHE	3.6
3	L	94	TRP	3.6
1	A	137	GLY	3.6
2	H	133	THR	3.6
1	A	118	PHE	3.6
2	H	94	TYR	3.6
1	A	106	LEU	3.6
1	A	363	VAL	3.6
1	A	362	ILE	3.5
1	A	406	VAL	3.5
1	A	353	ARG	3.5
1	A	101	ARG	3.5
1	A	304	LEU	3.5
1	A	176	GLU	3.5
1	A	311	ALA	3.5
1	A	264	VAL	3.5
1	A	43	PHE	3.5
1	A	188	VAL	3.4
3	L	32	ASN	3.4
1	A	262	VAL	3.4
3	L	143	ASP	3.4
1	A	306	PRO	3.4
1	A	386	PRO	3.4
1	A	183	VAL	3.4
1	A	263	ASP	3.4
1	A	351	LEU	3.4
1	A	258	ARG	3.3
3	L	50	TYR	3.3
1	A	260	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
3	L	185	GLU	3.3
1	A	76	TRP	3.3
1	A	174	PRO	3.2
1	A	330	VAL	3.2
1	A	111	TRP	3.2
1	A	47	ALA	3.2
1	A	163	TRP	3.1
2	H	49	VAL	3.1
1	A	298	THR	3.1
1	A	400	ASP	3.1
1	A	117	TRP	3.1
1	A	234	ILE	3.1
1	A	135	PHE	3.1
1	A	333	MET	3.1
1	A	270	LEU	3.1
1	A	116	VAL	3.1
1	A	202	ALA	3.1
1	A	294	ALA	3.1
1	A	199	PRO	3.1
2	H	53	SER	3.1
1	A	305	THR	3.1
1	A	412	THR	3.1
3	L	187	GLU	3.0
2	H	100	PRO	3.0
1	A	288	GLY	3.0
1	A	395	PHE	3.0
3	L	41	HIS	3.0
1	A	435	GLY	3.0
2	H	30	THR	3.0
3	L	111	ALA	3.0
1	A	134	GLU	3.0
1	A	148	LEU	2.9
1	A	99	SER	2.9
1	A	338	ILE	2.9
1	A	185	MET	2.9
2	H	2	VAL	2.9
2	H	37	VAL	2.9
1	A	60	PHE	2.9
3	L	80	THR	2.8
1	A	291	VAL	2.8
1	A	336	ASP	2.8
3	L	35	TRP	2.8

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Mol	Chain	Res	Type	RSRZ
3	L	183	LYS	2.8
1	A	413	VAL	2.8
1	A	420	VAL	2.8
1	A	182	PHE	2.8
3	L	36	TYR	2.8
1	A	396	THR	2.8
1	A	177	LEU	2.8
1	A	150	MET	2.7
1	A	181	ALA	2.7
1	A	172	THR	2.7
1	A	155	VAL	2.7
1	A	213	GLN	2.7
1	A	158	GLY	2.7
3	L	140	TYR	2.7
1	A	372	SER	2.7
3	L	191	SER	2.7
1	A	266	ALA	2.7
1	A	55	MET	2.7
1	A	164	LEU	2.7
1	A	206	ALA	2.7
2	H	174	GLY	2.7
3	L	150	ILE	2.7
3	L	40	SER	2.7
1	A	156	VAL	2.6
2	H	99	PRO	2.6
1	A	361	PHE	2.6
1	A	337	VAL	2.6
1	A	340	GLN	2.6
1	A	411	PHE	2.6
1	A	428	TYR	2.6
2	H	64	LYS	2.6
1	A	434	LEU	2.6
1	A	193	GLN	2.6
1	A	329	ALA	2.6
1	A	290	ASN	2.5
3	L	20	SER	2.5
1	A	390	VAL	2.5
1	A	278	MET	2.5
1	A	275	ASN	2.5
3	L	151	ASP	2.5
1	A	173	ARG	2.5
1	A	285	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	230	MET	2.5
3	L	195	GLU	2.5
1	A	97	ILE	2.5
1	A	214	GLY	2.4
1	A	417	VAL	2.4
1	A	184	LEU	2.4
2	H	128	GLY	2.4
1	A	153	GLY	2.4
3	L	28	SER	2.4
1	A	159	VAL	2.4
1	A	192	VAL	2.4
2	H	89	ASP	2.4
1	A	170	ARG	2.4
2	H	44	GLY	2.4
2	H	15	SER	2.4
1	A	175	LEU	2.4
3	L	31	ASN	2.4
1	A	355	ILE	2.4
2	H	212	GLU	2.4
2	H	87	THR	2.3
2	H	88	ASP	2.3
1	A	44	LEU	2.3
1	A	96	GLY	2.3
3	L	171	SER	2.3
1	A	354	LEU	2.3
1	A	309	GLY	2.3
3	L	189	HIS	2.3
1	A	394	LEU	2.3
3	L	10	THR	2.3
1	A	271	ALA	2.3
2	H	86	GLN	2.3
1	A	364	ILE	2.2
1	A	316	PHE	2.2
1	A	319	ALA	2.2
2	H	65	SER	2.2
1	A	57	PRO	2.2
2	H	132	THR	2.2
3	L	182	THR	2.2
3	L	30	SER	2.2
1	A	149	PRO	2.2
2	H	35	HIS	2.2
1	A	197	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
2	H	17	SER	2.2
1	A	359	PRO	2.2
1	A	98	ALA	2.1
1	A	273	LEU	2.2
1	A	141	ALA	2.1
1	A	431	TRP	2.1
2	H	52	TRP	2.1
1	A	391	PRO	2.1
3	L	29	ILE	2.1
3	L	51	VAL	2.1
1	A	56	ASP	2.1
3	L	14	THR	2.1
1	A	265	ILE	2.1
1	A	179	VAL	2.1
1	A	45	GLY	2.1
2	H	33	GLY	2.1
3	L	141	PRO	2.1
3	L	123	GLU	2.1
3	L	213	GLU	2.1
1	A	59	ASN	2.1
1	A	432	GLU	2.0
2	H	114	SER	2.0
3	L	156	GLN	2.0
3	L	166	GLN	2.0
1	A	104	PRO	2.0
3	L	165	ASP	2.0
1	A	51	SER	2.0
1	A	366	LEU	2.0
2	H	50	VAL	2.0
3	L	19	VAL	2.0
3	L	138	ASN	2.0
1	A	287	HIS	2.0
2	H	46	GLU	2.0
2	H	76	SER	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 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($\text{\AA}^2$)	Q<0.9
4	OS	A	502	1/1	0.88	0.10	184,184,184,184	0
4	OS	H	301	1/1	0.92	0.26	215,215,215,215	0
4	OS	A	501	1/1	0.95	0.14	190,190,190,190	0

6.5 Other polymers [i](#)

There are no such residues in this entry.