



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:37 PM UTC

PDB ID : 6XL1 / pdb_00006xl1
Title : crystal structure of cA4-activated Card1(D294N)
Authors : Rostol, J.; Xie, W.; Patel, D.J.; Marraffini, L.
Deposited on : 2020-06-27
Resolution : 1.95 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

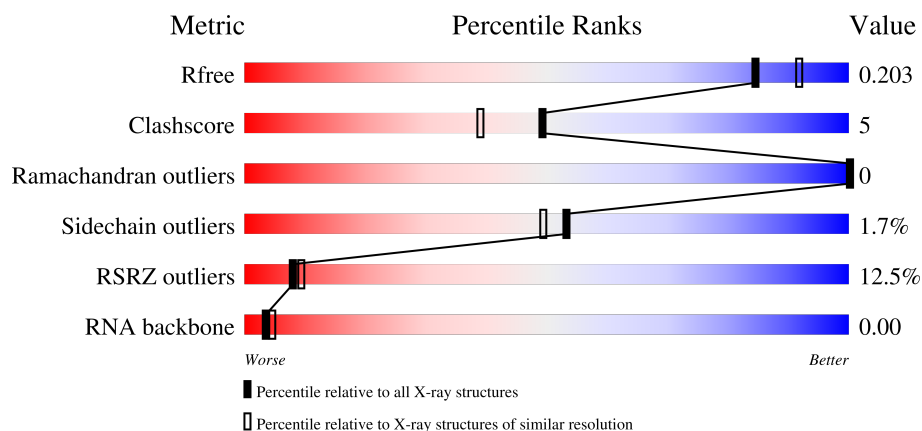
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)
RNA backbone	3983	1010 (2.32-1.60)

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	382	10% 87% 11% .
1	B	382	14% 84% 13% ..
2	C	4	25% 75%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6859 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 Card1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	373	Total	C	N	O	S	0	0	0
			3100	1996	508	585	11			
1	B	373	Total	C	N	O	S	0	0	0
			3100	1996	508	585	11			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	294	ASN	ASP	engineered mutation	UNP F2NWD3
A	374	GLY	-	expression tag	UNP F2NWD3
A	375	SER	-	expression tag	UNP F2NWD3
A	376	GLY	-	expression tag	UNP F2NWD3
A	377	HIS	-	expression tag	UNP F2NWD3
A	378	HIS	-	expression tag	UNP F2NWD3
A	379	HIS	-	expression tag	UNP F2NWD3
A	380	HIS	-	expression tag	UNP F2NWD3
A	381	HIS	-	expression tag	UNP F2NWD3
A	382	HIS	-	expression tag	UNP F2NWD3
B	294	ASN	ASP	engineered mutation	UNP F2NWD3
B	374	GLY	-	expression tag	UNP F2NWD3
B	375	SER	-	expression tag	UNP F2NWD3
B	376	GLY	-	expression tag	UNP F2NWD3
B	377	HIS	-	expression tag	UNP F2NWD3
B	378	HIS	-	expression tag	UNP F2NWD3
B	379	HIS	-	expression tag	UNP F2NWD3
B	380	HIS	-	expression tag	UNP F2NWD3
B	381	HIS	-	expression tag	UNP F2NWD3
B	382	HIS	-	expression tag	UNP F2NWD3

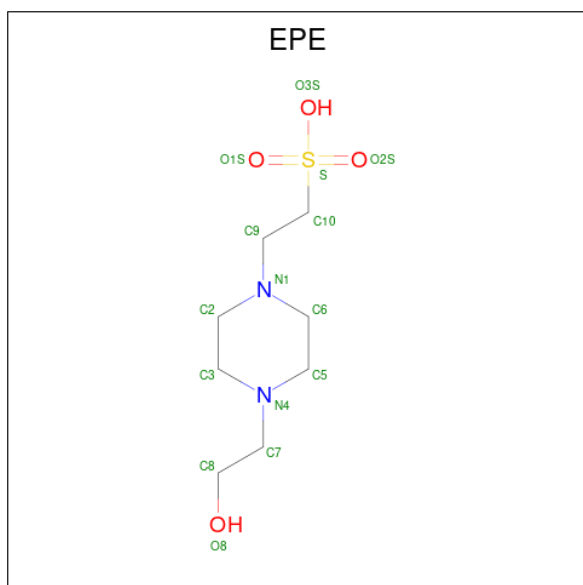
- Molecule 2 is a RNA chain called cA4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	P	0	0	0
			88	40	20	24	4			

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	B	3	Total	Mn	0	0
			3	3		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	265	Total	O	0	0
			265	265		
5	B	262	Total	O	0	0
			262	262		

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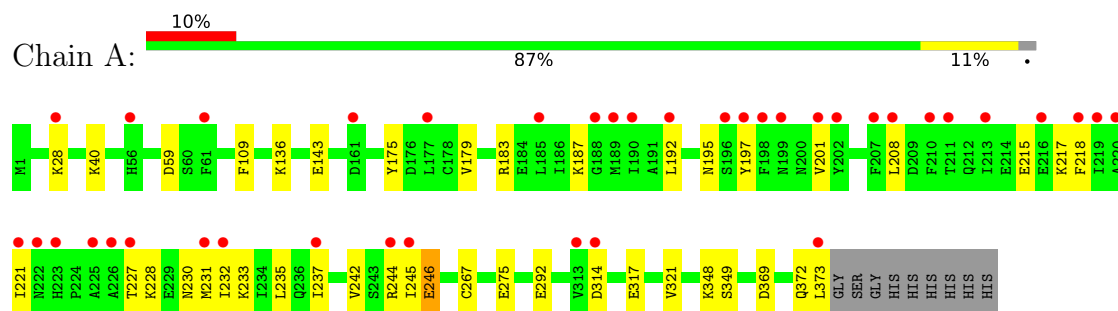
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	25	Total	O	0	0
			25	25		

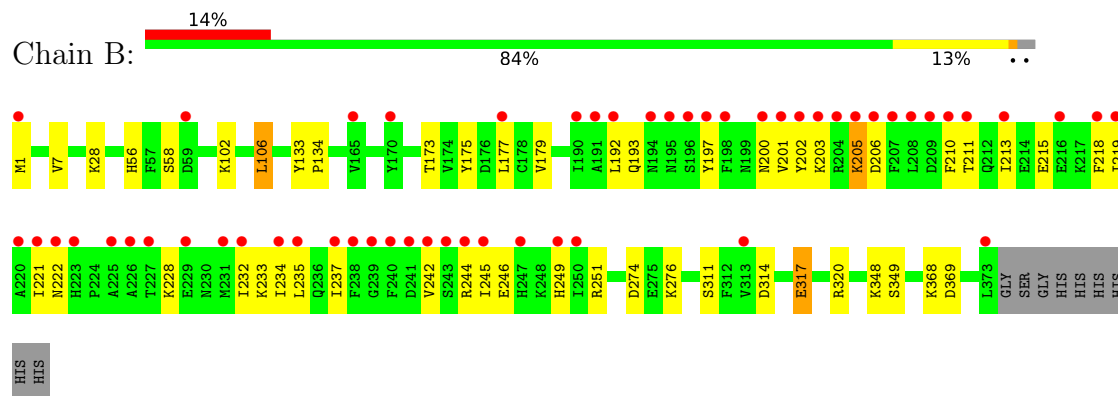
3 Residue-property plots

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: Card1



• Molecule 1: Card1



• Molecule 2: cA4



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	111.95Å 111.95Å 172.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.58 – 1.95 39.58 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.58-1.95) 99.9 (39.58-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.18_3855, PHENIX 1.18_3855	Depositor
R, R_{free}	0.178 , 0.202 0.179 , 0.203	Depositor DCC
R_{free} test set	4045 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6859	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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: EPE, 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	0.31	0/3165	0.48	0/4267
1	B	0.38	0/3165	0.48	0/4267
2	C	0.72	0/99	0.84	0/152
All	All	0.36	0/6429	0.49	0/8686

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	3100	0	3077	34	0
1	B	3100	0	3077	34	0
2	C	88	0	44	0	0
3	A	1	0	0	0	0
3	B	3	0	0	0	0
4	B	15	0	18	2	0
5	A	265	0	0	8	3
5	B	262	0	0	2	3
5	C	25	0	0	0	0
All	All	6859	0	6216	67	3

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 (67) 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:195:ASN:HD22	1:A:221:ILE:HA	1.33	0.93
1:B:368:LYS:HE2	4:B:401:EPE:H52	1.71	0.73
1:A:143:GLU:OE2	5:A:1001:HOH:O	2.06	0.71
1:B:173:THR:OG1	5:B:501:HOH:O	2.11	0.68
1:A:195:ASN:ND2	1:A:221:ILE:HA	2.09	0.65
1:A:197:TYR:CD1	1:A:217:LYS:HG2	2.33	0.64
1:B:232:ILE:HG23	1:B:242:VAL:HG21	1.80	0.64
1:B:246:GLU:HG3	1:B:249:HIS:CE1	2.33	0.63
1:A:201:VAL:HG11	1:A:208:LEU:HD11	1.82	0.61
1:B:202:TYR:OH	1:B:245:ILE:O	2.15	0.59
1:A:235:LEU:HD22	1:A:245:ILE:HD12	1.84	0.58
1:A:183:ARG:NH2	5:A:1003:HOH:O	2.34	0.57
1:B:56:HIS:O	1:B:56:HIS:ND1	2.38	0.57
1:A:59:ASP:OD1	5:A:1002:HOH:O	2.18	0.56
1:B:233:LYS:O	1:B:237:ILE:HG13	2.05	0.56
1:B:102:LYS:O	1:B:106:LEU:HB2	2.04	0.56
1:B:276:LYS:NZ	5:B:502:HOH:O	2.21	0.55
1:B:210:PHE:O	1:B:213:ILE:HG22	2.07	0.55
1:A:183:ARG:O	1:A:187:LYS:HG2	2.08	0.54
1:A:227:THR:OG1	1:A:230:ASN:OD1	2.26	0.53
1:B:210:PHE:HB2	1:B:242:VAL:HG13	1.90	0.53
1:A:317:GLU:O	1:A:321:VAL:HG23	2.09	0.52
1:A:215:GLU:OE1	1:A:215:GLU:N	2.38	0.52
1:A:109:PHE:CD2	1:B:106:LEU:HD21	2.46	0.50
1:B:317:GLU:OE1	1:B:320:ARG:NH2	2.46	0.49
1:A:232:ILE:HG23	1:A:242:VAL:HG11	1.94	0.48
1:B:197:TYR:O	1:B:201:VAL:HG12	2.14	0.48
1:A:28:LYS:NZ	5:A:1016:HOH:O	2.46	0.48
1:B:232:ILE:HA	1:B:235:LEU:HD12	1.95	0.48
1:A:192:LEU:HD12	1:A:231:MET:HG3	1.96	0.48
1:A:267:CYS:HB3	1:A:275:GLU:HG3	1.96	0.48
1:B:200:ASN:HA	1:B:203:LYS:HE2	1.96	0.47
1:A:314:ASP:OD2	5:A:1003:HOH:O	2.20	0.47
1:B:173:THR:HB	1:B:177:LEU:HD13	1.97	0.47
1:B:215:GLU:O	1:B:219:ILE:HD12	2.15	0.46
1:A:175:TYR:HA	1:A:179:VAL:HB	1.96	0.46
1:B:274:ASP:OD1	1:B:276:LYS:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:TYR:CZ	1:A:217:LYS:HE2	2.51	0.45
1:B:205:LYS:HE3	1:B:206:ASP:H	1.81	0.45
1:B:175:TYR:HA	1:B:179:VAL:HB	1.99	0.45
1:A:197:TYR:CE2	1:A:217:LYS:HE2	2.52	0.44
1:B:246:GLU:HG3	1:B:249:HIS:HE1	1.79	0.44
1:A:192:LEU:HG	1:A:221:ILE:HD11	2.00	0.43
1:B:222:ASN:C	1:B:222:ASN:HD22	2.26	0.43
1:B:211:THR:O	1:B:228:LYS:NZ	2.51	0.43
1:A:348:LYS:NZ	1:A:373:LEU:O	2.47	0.43
1:A:244:ARG:NH2	1:A:246:GLU:OE2	2.40	0.43
1:A:242:VAL:HG22	5:A:1050:HOH:O	2.19	0.42
1:A:348:LYS:HA	1:A:348:LYS:HD2	1.84	0.42
1:A:215:GLU:HG3	1:A:228:LYS:HD3	2.02	0.42
1:A:218:PHE:CE1	1:A:228:LYS:HA	2.54	0.42
1:A:231:MET:HA	1:A:231:MET:HE2	2.01	0.42
1:B:218:PHE:CE2	1:B:228:LYS:HG3	2.55	0.42
1:A:218:PHE:CD1	1:A:228:LYS:HA	2.55	0.42
1:A:136:LYS:HE3	5:A:1083:HOH:O	2.20	0.42
1:B:349:SER:O	1:B:369:ASP:HB2	2.19	0.42
1:B:193:GLN:HE22	1:B:251:ARG:HG2	1.85	0.42
1:A:349:SER:O	1:A:369:ASP:HB2	2.20	0.42
1:B:28:LYS:HB2	1:B:28:LYS:HE2	1.89	0.41
1:B:200:ASN:HA	1:B:203:LYS:CE	2.51	0.41
1:B:192:LEU:HA	1:B:221:ILE:HD13	2.02	0.41
1:A:292:GLU:HB3	5:A:1019:HOH:O	2.21	0.41
1:B:133:TYR:CD2	1:B:134:PRO:HA	2.56	0.41
1:B:234:ILE:HD12	1:B:234:ILE:HA	1.89	0.41
1:B:348:LYS:HA	1:B:348:LYS:HD2	1.89	0.40
1:A:233:LYS:O	1:A:237:ILE:HG13	2.21	0.40
1:B:368:LYS:HE3	4:B:401:EPE:H82	2.04	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1205:HOH:O	5:B:502:HOH:O[6_554]	1.96	0.24
5:A:1191:HOH:O	5:B:578:HOH:O[3_555]	2.04	0.16
5:A:1001:HOH:O	5:B:728:HOH:O[6_554]	2.17	0.03

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	371/382 (97%)	363 (98%)	8 (2%)	0	100	100
1	B	371/382 (97%)	362 (98%)	9 (2%)	0	100	100
All	All	742/764 (97%)	725 (98%)	17 (2%)	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	345/352 (98%)	342 (99%)	3 (1%)	70	70
1	B	345/352 (98%)	336 (97%)	9 (3%)	40	33
All	All	690/704 (98%)	678 (98%)	12 (2%)	53	49

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LYS
1	A	246	GLU
1	A	372	GLN
1	B	1	MET
1	B	7	VAL
1	B	58	SER
1	B	106	LEU
1	B	205	LYS

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Mol	Chain	Res	Type
1	B	244	ARG
1	B	311	SER
1	B	314	ASP
1	B	317	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	ASN
1	A	212	GLN
1	A	281	ASN
1	A	291	ASN
1	B	114	ASN
1	B	137	GLN
1	B	194	ASN
1	B	195	ASN
1	B	200	ASN
1	B	222	ASN
1	B	236	GLN
1	B	281	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	3/4 (75%)	3 (100%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	2	A
2	C	3	A
2	C	4	A

There are no RNA pucker outliers to report.

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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
4	EPE	B	401	-	15,15,15	1.39	2 (13%)	19,20,20	2.00	5 (26%)

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
4	EPE	B	401	-	-	4/9/19/19	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	401	EPE	C10-S	3.60	1.82	1.77
4	B	401	EPE	O1S-S	2.50	1.52	1.45

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	EPE	C6-C5-N4	4.20	119.11	110.65
4	B	401	EPE	C5-N4-C3	4.16	117.79	108.84
4	B	401	EPE	O1S-S-C10	3.57	112.12	106.73
4	B	401	EPE	O3S-S-O1S	-3.31	103.12	111.40
4	B	401	EPE	O2S-S-C10	2.63	110.70	106.73

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	401	EPE	C9-C10-S-O2S
4	B	401	EPE	C9-C10-S-O3S
4	B	401	EPE	C10-C9-N1-C2
4	B	401	EPE	C9-C10-S-O1S

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	401	EPE	2	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

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	373/382 (97%)	0.37	39 (10%) 11 13	21, 33, 72, 83	0
1	B	373/382 (97%)	0.43	55 (14%) 6 6	22, 33, 77, 84	0
2	C	4/4 (100%)	-1.16	0 100 100	22, 22, 23, 24	0
All	All	750/768 (97%)	0.39	94 (12%) 8 9	21, 33, 75, 84	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	208	LEU	5.7
1	B	210	PHE	5.2
1	B	213	ILE	4.6
1	B	250	ILE	4.6
1	B	240	PHE	4.3
1	B	245	ILE	4.1
1	A	213	ILE	4.0
1	B	201	VAL	4.0
1	B	194	ASN	3.9
1	A	197	TYR	3.9
1	B	239	GLY	3.8
1	B	202	TYR	3.8
1	A	218	PHE	3.8
1	B	235	LEU	3.6
1	B	204	ARG	3.6
1	A	185	LEU	3.5
1	A	219	ILE	3.4
1	A	201	VAL	3.4
1	A	313	VAL	3.4
1	B	200	ASN	3.4
1	B	207	PHE	3.4
1	A	56	HIS	3.3
1	B	242	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	197	TYR	3.3
1	A	192	LEU	3.3
1	A	196	SER	3.3
1	A	221	ILE	3.2
1	B	198	PHE	3.2
1	B	313	VAL	3.2
1	B	218	PHE	3.2
1	B	232	ILE	3.2
1	A	211	THR	3.2
1	A	220	ALA	3.1
1	B	234	ILE	3.1
1	B	211	THR	3.0
1	B	196	SER	2.9
1	A	244	ARG	2.9
1	A	208	LEU	2.9
1	B	192	LEU	2.9
1	B	1	MET	2.9
1	A	28	LYS	2.9
1	A	210	PHE	2.8
1	A	199	ASN	2.8
1	A	207	PHE	2.8
1	A	189	MET	2.8
1	B	225	ALA	2.8
1	B	219	ILE	2.7
1	B	221	ILE	2.7
1	A	216	GLU	2.7
1	A	198	PHE	2.7
1	B	177	LEU	2.7
1	B	244	ARG	2.7
1	B	209	ASP	2.6
1	A	177	LEU	2.6
1	B	249	HIS	2.6
1	B	231	MET	2.6
1	A	225	ALA	2.6
1	B	170	TYR	2.5
1	A	232	ILE	2.5
1	B	205	LYS	2.4
1	B	227	THR	2.4
1	A	373	LEU	2.4
1	B	203	LYS	2.4
1	A	190	ILE	2.4
1	B	165	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	245	ILE	2.3
1	B	226	ALA	2.3
1	B	220	ALA	2.3
1	B	216	GLU	2.3
1	B	195	ASN	2.3
1	A	231	MET	2.3
1	A	202	TYR	2.3
1	A	237	ILE	2.3
1	B	206	ASP	2.3
1	A	61	PHE	2.2
1	B	237	ILE	2.2
1	A	188	GLY	2.2
1	B	190	ILE	2.2
1	A	227	THR	2.2
1	A	161	ASP	2.2
1	A	226	ALA	2.2
1	B	373	LEU	2.1
1	A	223	HIS	2.1
1	A	314	ASP	2.1
1	B	241	ASP	2.1
1	B	247	HIS	2.1
1	A	222	ASN	2.1
1	B	238	PHE	2.1
1	B	243	SER	2.1
1	B	222	ASN	2.1
1	B	223	HIS	2.0
1	B	191	ALA	2.0
1	B	59	ASP	2.0
1	B	229	GLU	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands

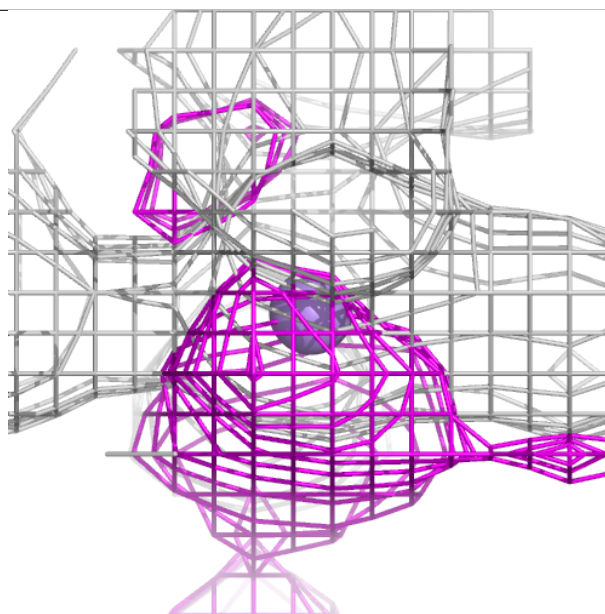
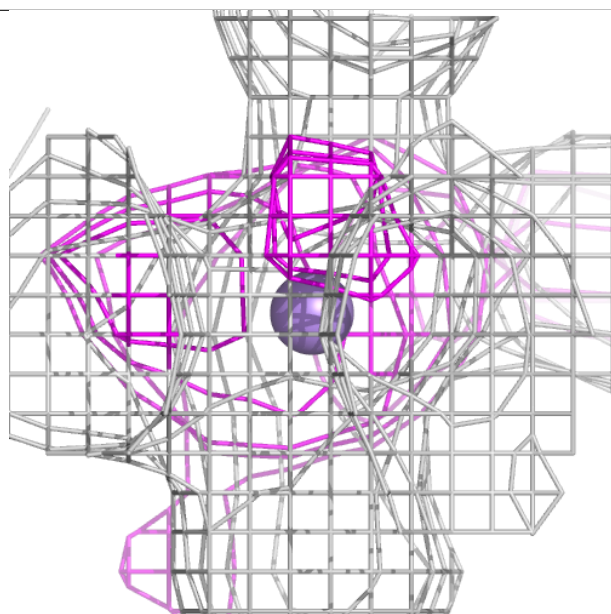
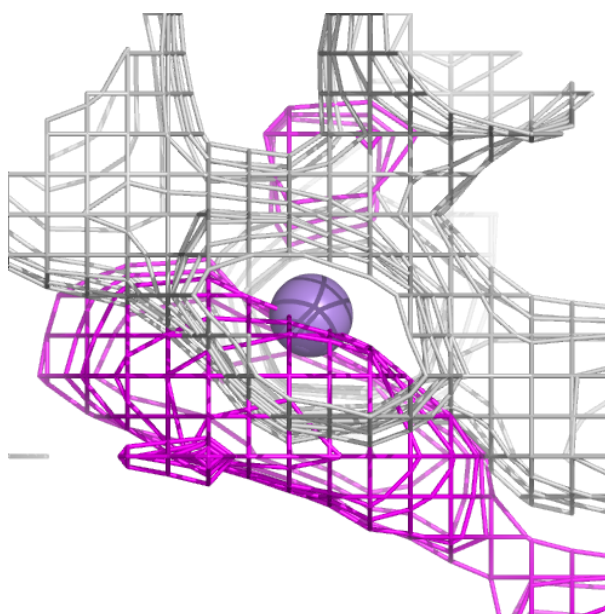
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EPE	B	401	15/15	0.80	0.19	41,54,60,63	0
3	MN	A	901	1/1	0.88	0.26	81,81,81,81	0
3	MN	B	402	1/1	0.95	0.27	66,66,66,66	0
3	MN	B	404	1/1	0.99	0.06	43,43,43,43	0
3	MN	B	403	1/1	1.00	0.02	30,30,30,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

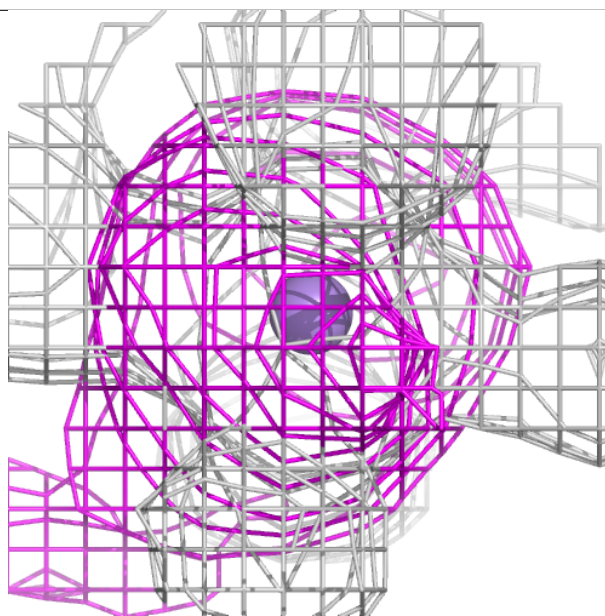
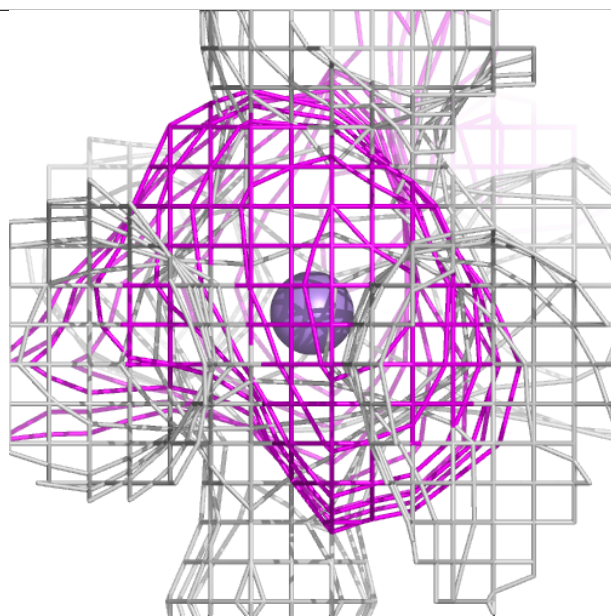
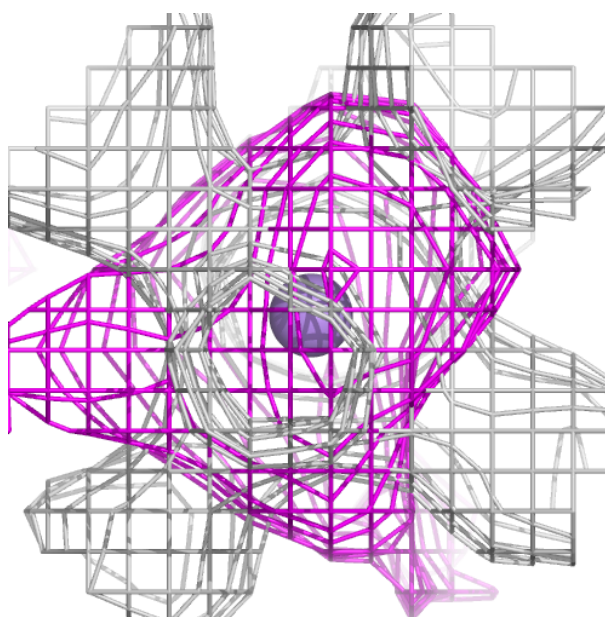
Electron density around MN A 901:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



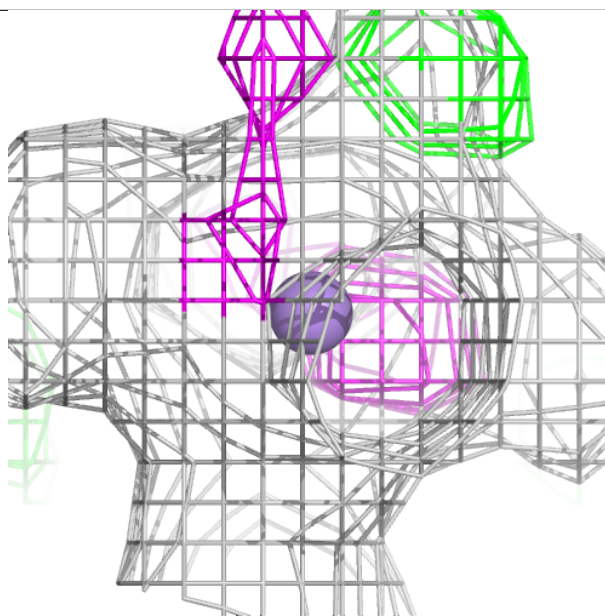
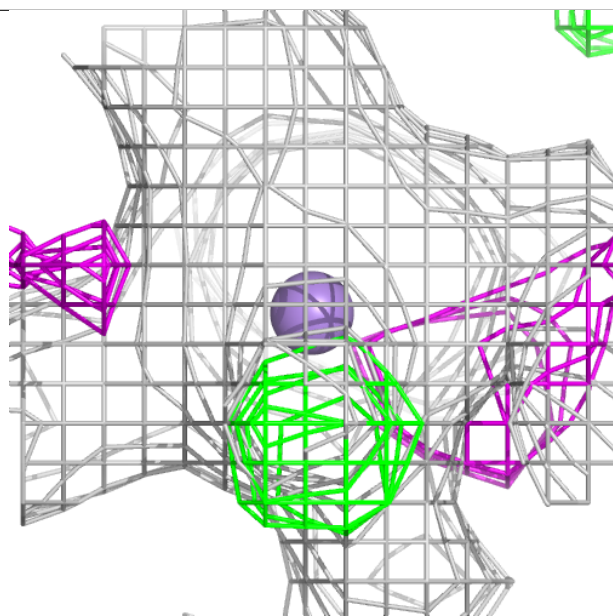
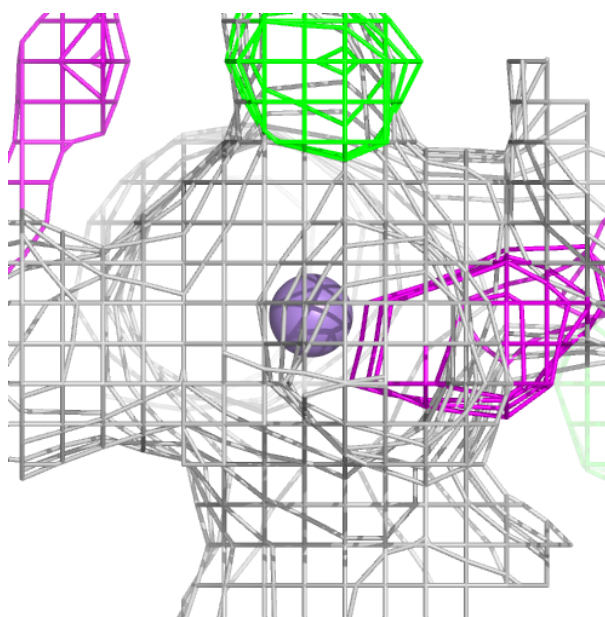
Electron density around MN B 402:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



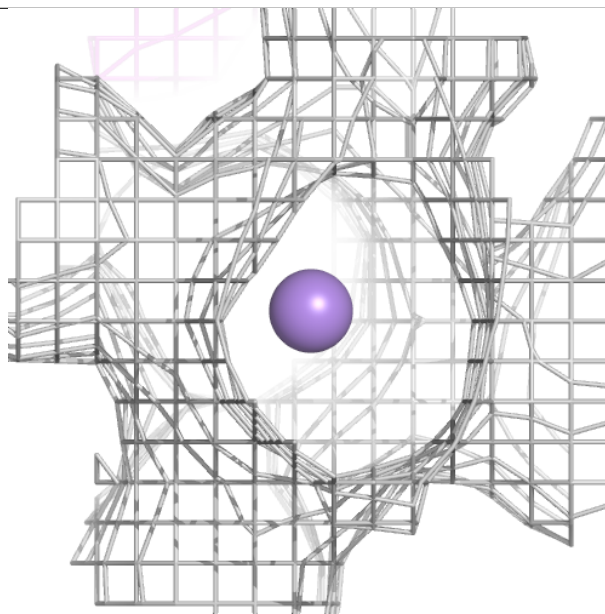
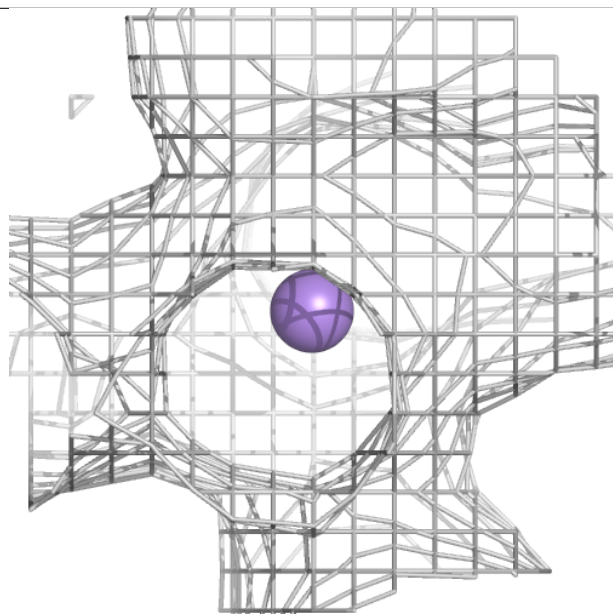
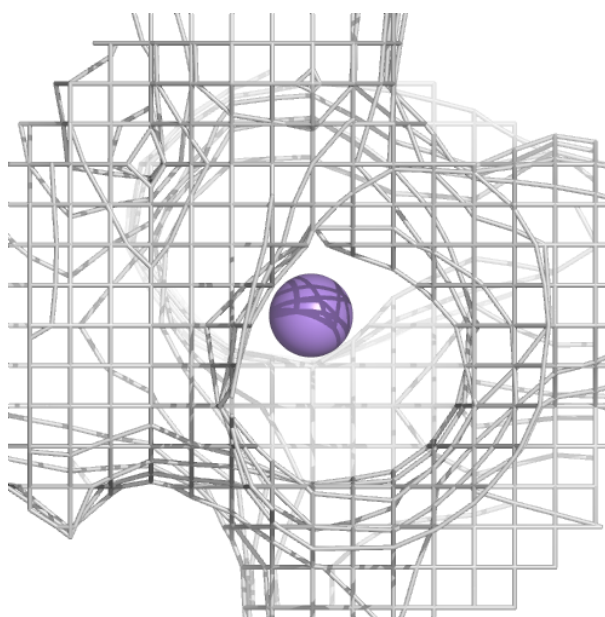
Electron density around MN B 404:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MN B 403:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.