



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:15 PM UTC

PDB ID : 8D8W / pdb_00008d8w
Title : Crystal structure of ChoE with Ser38 adopting alternative conformations
Authors : Pham, V.D.; Shi, R.
Deposited on : 2022-06-09
Resolution : 1.40 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

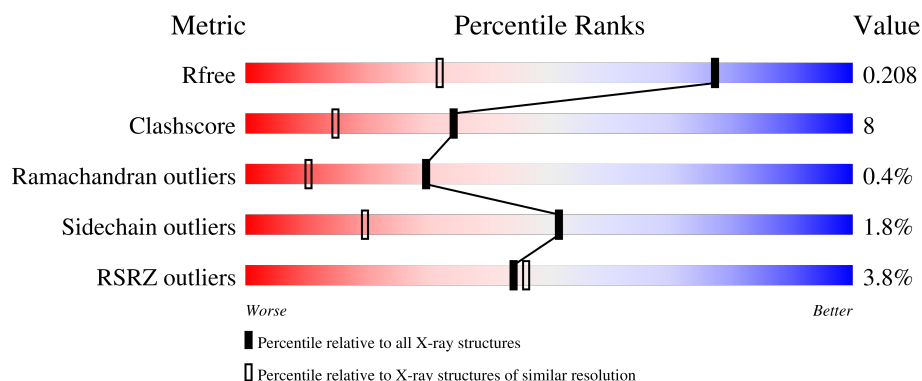
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	2563 (1.40-1.40)
Clashscore	190562	2660 (1.40-1.40)
Ramachandran outliers	187476	2611 (1.40-1.40)
Sidechain outliers	187428	2610 (1.40-1.40)
RSRZ outliers	180081	2561 (1.40-1.40)

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	287	6% 87% 10%
1	B	287	2% 85% 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IOD	A	406	-	-	X	-
3	IOD	B	406	-	-	X	-

2 Entry composition [i](#)

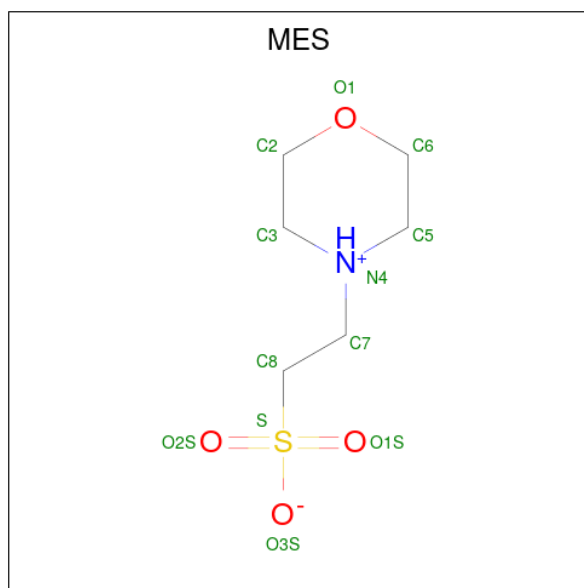
There are 4 unique types of molecules in this entry. The entry contains 5245 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ChoE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	6	0
			2257	1417	422	413	5			
1	B	287	Total	C	N	O	S	0	6	0
			2256	1420	419	412	5			

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



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

- Molecule 3 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total 6	I 6	0	2
3	B	8	Total 8	I 8	0	2

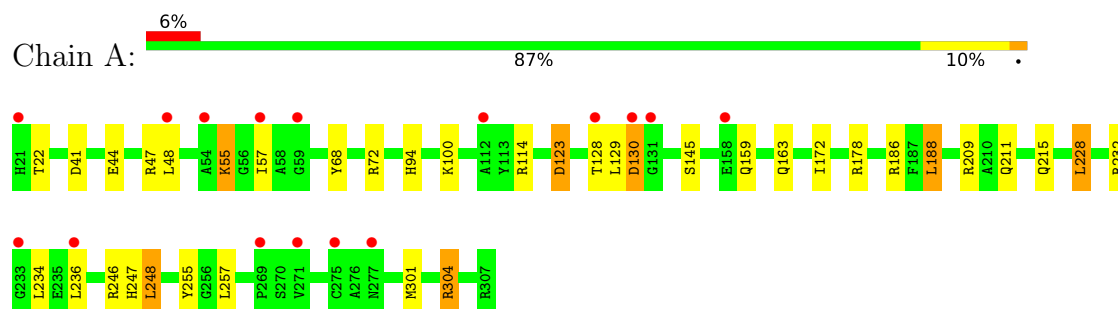
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	296	Total 296	O 296	0	0
4	B	398	Total 398	O 398	0	0

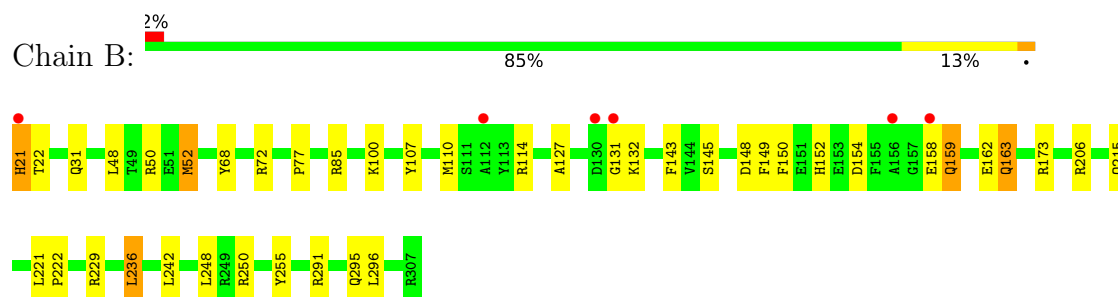
3 Residue-property plots [i](#)

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: ChoE



• Molecule 1: ChoE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	45.72Å 81.57Å 81.49Å 90.00° 99.89° 90.00°	Depositor
Resolution (Å)	45.08 – 1.40 45.08 – 1.40	Depositor EDS
% Data completeness (in resolution range)	94.8 (45.08-1.40) 94.8 (45.08-1.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.169 , 0.197 (Not available) , 0.208	Depositor DCC
R_{free} test set	5502 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	11.7	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.3	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	5245	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.10% 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: IOD, MES

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	1.26	2/2332 (0.1%)	1.43	12/3170 (0.4%)
1	B	1.33	9/2331 (0.4%)	1.41	10/3169 (0.3%)
All	All	1.29	11/4663 (0.2%)	1.42	22/6339 (0.3%)

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
1	A	0	1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	130	ASP	C-O	-7.07	1.15	1.24
1	B	206	ARG	NE-CZ	-6.60	1.25	1.33
1	B	152	HIS	C-O	6.52	1.31	1.24
1	B	236	LEU	C-O	6.50	1.31	1.23
1	B	152	HIS	ND1-CE1	6.27	1.38	1.32
1	B	127	ALA	C-O	6.08	1.32	1.24
1	A	94	HIS	CE1-NE2	5.86	1.38	1.32
1	B	149	PHE	C-O	-5.38	1.17	1.24
1	B	31	GLN	C-O	-5.33	1.18	1.23
1	B	154	ASP	C-O	5.19	1.30	1.24
1	B	21	HIS	CE1-NE2	5.15	1.37	1.32

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	291	ARG	CG-CD-NE	-7.95	94.52	112.00
1	B	173	ARG	NE-CZ-NH2	7.78	126.20	119.20
1	B	229	ARG	NE-CZ-NH2	7.28	125.75	119.20
1	B	52	MET	CG-SD-CE	6.90	116.08	100.90
1	A	186	ARG	NE-CZ-NH1	-6.79	114.71	121.50
1	B	173	ARG	NE-CZ-NH1	-6.48	115.02	121.50
1	A	123	ASP	CA-CB-CG	5.90	118.50	112.60
1	A	211	GLN	CB-CG-CD	5.72	122.32	112.60
1	B	229	ARG	NE-CZ-NH1	-5.69	115.81	121.50
1	A	22	THR	CA-CB-OG1	-5.66	101.11	109.60
1	B	148	ASP	CA-CB-CG	5.63	118.23	112.60
1	B	250	ARG	CA-C-N	5.60	127.99	122.28
1	B	250	ARG	C-N-CA	5.60	127.99	122.28
1	A	41	ASP	CA-CB-CG	5.48	118.08	112.60
1	A	232	ARG	CA-C-O	-5.40	113.03	119.35
1	A	209	ARG	NE-CZ-NH2	5.36	124.02	119.20
1	A	129	LEU	CA-C-N	-5.30	111.42	121.54
1	A	129	LEU	C-N-CA	-5.30	111.42	121.54
1	B	143	PHE	CA-CB-CG	5.23	119.03	113.80
1	A	172	ILE	N-CA-C	-5.12	105.72	110.53
1	A	246	ARG	N-CA-C	-5.04	105.79	111.28
1	A	128	THR	CA-CB-OG1	-5.02	102.07	109.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	130	ASP	Peptide

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	2257	0	2195	25	0
1	B	2256	0	2199	37	0
2	A	12	0	13	2	0
2	B	12	0	12	3	0
3	A	6	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	0	4	0
4	A	296	0	0	11	0
4	B	398	0	0	14	0
All	All	5245	0	4419	68	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 8.

All (68) 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:B:48[B]:LEU:HG	1:B:52:MET:HE3	1.28	1.07
1:B:48[B]:LEU:CG	1:B:52:MET:HE3	1.89	1.03
1:A:228:LEU:HD22	4:A:724:HOH:O	1.58	1.02
1:B:114:ARG:NE	3:B:406:IOD:I	2.65	0.99
1:A:114:ARG:NE	3:A:406:IOD:I	2.74	0.91
3:B:403[A]:IOD:I	4:B:734:HOH:O	2.61	0.88
1:B:131:GLY:C	4:B:505:HOH:O	2.18	0.85
1:B:50:ARG:NH1	4:B:501:HOH:O	2.10	0.82
3:A:403[A]:IOD:I	4:A:650:HOH:O	2.75	0.74
2:B:401:MES:H32	4:B:791:HOH:O	1.88	0.74
1:B:48[B]:LEU:CD2	1:B:52:MET:HE3	2.20	0.71
3:A:406:IOD:I	4:A:505:HOH:O	2.83	0.66
1:A:123:ASP:OD1	1:A:178:ARG:NH2	2.27	0.66
1:A:48:LEU:HD11	4:A:505:HOH:O	1.95	0.66
1:B:162:GLU:HG3	4:B:750:HOH:O	1.99	0.62
1:A:248:LEU:HD22	1:A:257:LEU:HD12	1.80	0.62
1:A:44:GLU:OE1	1:A:47:ARG:NH2	2.32	0.61
1:B:21:HIS:HD2	1:B:22:THR:O	1.82	0.61
1:B:100:LYS:HE3	3:B:404[B]:IOD:I	2.71	0.60
1:A:188:LEU:HD11	1:A:304[A]:ARG:NH2	2.16	0.60
1:B:132:LYS:N	4:B:505:HOH:O	2.29	0.59
1:B:48[B]:LEU:HD21	1:B:52:MET:CE	2.34	0.58
1:B:248[A]:LEU:HD23	1:B:255:TYR:CD2	2.39	0.57
1:A:234:LEU:HD11	4:A:724:HOH:O	2.06	0.56
1:B:159:GLN:HB2	4:B:503:HOH:O	2.05	0.55
1:B:159:GLN:NE2	4:B:503:HOH:O	2.22	0.55
2:A:401:MES:H32	4:A:742:HOH:O	2.08	0.54
2:A:401:MES:C6	4:A:507:HOH:O	2.57	0.53
1:A:159:GLN:NE2	4:A:502:HOH:O	2.30	0.52
1:A:114:ARG:NE	4:A:505:HOH:O	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85[A]:ARG:HH22	1:B:295:GLN:HE22	1.58	0.52
1:A:68:TYR:HB3	1:A:72:ARG:O	2.11	0.50
1:B:48[B]:LEU:HD21	1:B:52:MET:HE3	1.94	0.50
1:B:77:PRO:CB	1:B:85[B]:ARG:HH12	2.25	0.48
1:A:248:LEU:HD23	1:A:255:TYR:CD2	2.49	0.47
1:B:248[B]:LEU:CD1	4:B:821:HOH:O	2.62	0.47
1:A:236:LEU:C	1:A:236:LEU:HD23	2.41	0.46
1:A:159:GLN:HB2	4:A:502:HOH:O	2.16	0.45
1:B:114:ARG:CZ	3:B:406:IOD:I	3.34	0.45
1:B:242:LEU:CD2	4:B:694:HOH:O	2.63	0.45
1:B:68:TYR:HB3	1:B:72:ARG:O	2.17	0.45
1:B:159:GLN:HG2	1:B:163:GLN:HB3	1.99	0.44
1:B:85[A]:ARG:HH22	1:B:295:GLN:NE2	2.16	0.44
1:A:188:LEU:C	1:A:188:LEU:HD23	2.42	0.44
1:B:77:PRO:HB3	1:B:85[B]:ARG:HH12	1.82	0.43
1:A:100:LYS:HE3	3:A:404[B]:IOD:I	2.87	0.43
1:B:48[B]:LEU:CD2	1:B:52:MET:CE	2.89	0.43
1:B:150:PHE:CD2	2:B:401:MES:H61	2.53	0.43
1:A:247:HIS:NE2	1:A:255:TYR:OH	2.51	0.43
1:B:215:GLN:NE2	4:B:513:HOH:O	2.51	0.43
1:A:55:LYS:HB2	1:A:57:ILE:CD1	2.48	0.43
1:B:85[B]:ARG:NH1	4:B:511:HOH:O	2.50	0.43
1:A:301:MET:HA	1:A:304[A]:ARG:HH21	1.83	0.43
1:B:236:LEU:C	1:B:236:LEU:HD23	2.44	0.43
1:B:248[B]:LEU:HD11	1:B:296:LEU:HD13	2.01	0.42
1:B:242:LEU:HD23	4:B:694:HOH:O	2.19	0.42
1:B:159:GLN:HB2	1:B:159:GLN:HE21	1.56	0.42
1:A:159:GLN:HG2	1:A:163:GLN:HB3	2.01	0.42
1:B:221:LEU:N	1:B:222:PRO:HD2	2.35	0.42
1:B:52:MET:HE1	1:B:110:MET:HG2	2.02	0.41
1:B:85[A]:ARG:NH2	1:B:295:GLN:OE1	2.53	0.41
1:A:248:LEU:HD22	1:A:257:LEU:CD1	2.50	0.41
1:A:188:LEU:HD11	1:A:304[A]:ARG:HH22	1.84	0.41
1:A:301:MET:O	1:A:304[A]:ARG:NE	2.53	0.41
1:B:107:TYR:OH	2:B:401:MES:H31	2.21	0.40
1:A:304[B]:ARG:HD2	1:A:304[B]:ARG:HA	1.92	0.40
1:B:248[B]:LEU:HD13	4:B:821:HOH:O	2.20	0.40
1:A:215:GLN:NE2	4:A:517:HOH:O	2.54	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	A	291/287 (101%)	278 (96%)	12 (4%)	1 (0%)	36	16
1	B	291/287 (101%)	283 (97%)	7 (2%)	1 (0%)	36	16
All	All	582/574 (101%)	561 (96%)	19 (3%)	2 (0%)	30	16

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	SER
1	B	145	SER

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	227/221 (103%)	221 (97%)	6 (3%)	40	11
1	B	227/221 (103%)	224 (99%)	3 (1%)	61	30
All	All	454/442 (103%)	445 (98%)	9 (2%)	51	17

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	LYS
1	A	188	LEU
1	A	228	LEU
1	A	248	LEU

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Mol	Chain	Res	Type
1	A	304[A]	ARG
1	A	304[B]	ARG
1	B	158	GLU
1	B	159	GLN
1	B	163	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	61	GLN
1	A	163	GLN
1	A	215	GLN
1	A	268	GLN
1	B	21	HIS
1	B	61	GLN
1	B	215	GLN
1	B	268	GLN

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 16 ligands modelled in this entry, 14 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	MES	B	401	-	12,12,12	0.83	1 (8%)	15,16,16	1.85	5 (33%)
2	MES	A	401	-	12,12,12	0.91	0	15,16,16	1.20	2 (13%)

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	MES	B	401	-	-	3/6/14/14	0/1/1/1
2	MES	A	401	-	-	3/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	MES	C7-N4	2.01	1.52	1.47

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	MES	C6-C5-N4	-3.85	104.27	110.12
2	B	401	MES	C7-N4-C3	3.43	120.37	111.24
2	A	401	MES	C7-N4-C3	2.61	118.19	111.24
2	B	401	MES	O1-C2-C3	2.45	117.05	111.77
2	A	401	MES	O1-C6-C5	-2.13	107.19	111.77
2	B	401	MES	C2-C3-N4	2.06	113.26	110.12
2	B	401	MES	C7-N4-C5	2.05	116.70	111.24

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	MES	C8-C7-N4-C3
2	B	401	MES	C8-C7-N4-C3
2	A	401	MES	N4-C7-C8-S
2	B	401	MES	N4-C7-C8-S

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Mol	Chain	Res	Type	Atoms
2	A	401	MES	C8-C7-N4-C5
2	B	401	MES	C8-C7-N4-C5

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	MES	3	0
2	A	401	MES	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	287/287 (100%)	0.22	16 (5%) 30 32	7, 16, 36, 55	6 (2%)
1	B	287/287 (100%)	-0.23	6 (2%) 63 66	4, 11, 24, 46	6 (2%)
All	All	574/574 (100%)	-0.01	22 (3%) 44 46	4, 13, 32, 55	12 (2%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	130	ASP	4.2
1	A	112	ALA	3.6
1	A	158	GLU	3.1
1	A	271	VAL	3.1
1	B	131	GLY	2.9
1	A	57	ILE	2.8
1	B	158	GLU	2.7
1	A	233	GLY	2.5
1	A	48	LEU	2.5
1	A	236	LEU	2.4
1	A	131	GLY	2.4
1	A	21	HIS	2.3
1	B	112	ALA	2.3
1	B	130	ASP	2.2
1	B	21	HIS	2.2
1	B	156	ALA	2.2
1	A	269	PRO	2.2
1	A	128	THR	2.1
1	A	275	CYS	2.1
1	A	59	GLY	2.0
1	A	54	ALA	2.0
1	A	277	ASN	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($\text{\AA}^2$)	Q<0.9
2	MES	B	401	12/12	0.86	0.14	23,28,31,34	0
2	MES	A	401	12/12	0.87	0.14	26,34,38,40	0
3	IOD	B	408	1/1	0.90	0.16	44,44,44,44	1
3	IOD	B	404[B]	1/1	0.93	0.15	36,36,36,36	1
3	IOD	A	406	1/1	0.93	0.16	48,48,48,48	1
3	IOD	A	404[B]	1/1	0.94	0.14	40,40,40,40	1
3	IOD	A	407	1/1	0.94	0.17	48,48,48,48	1
3	IOD	A	405	1/1	0.96	0.08	40,40,40,40	1
3	IOD	B	405	1/1	0.97	0.09	31,31,31,31	1
3	IOD	B	407	1/1	0.97	0.16	39,39,39,39	1
3	IOD	A	403[A]	1/1	0.97	0.10	30,30,30,30	1
3	IOD	B	409	1/1	0.97	0.20	43,43,43,43	1
3	IOD	B	406	1/1	0.98	0.19	29,29,29,29	1
3	IOD	B	403[A]	1/1	0.98	0.10	22,22,22,22	1
3	IOD	B	402	1/1	0.99	0.04	19,19,19,19	1
3	IOD	A	402	1/1	0.99	0.06	27,27,27,27	1

6.5 Other polymers [i](#)

There are no such residues in this entry.