



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:05 PM UTC

PDB ID : 4N3P / pdb_00004n3p
Title : Crystal Structure of De Novo designed Serine Hydrolase OSH18, Northeast Structural Genomics Consortium (NESG) Target OR396
Authors : Kuzin, A.; Lew, S.; Rajagopalan, S.; Seetharaman, J.; Mao, L.; Xiao, R.; Kogan, S.; Maglaqui, M.; Everett, J.K.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2013-10-07
Resolution : 2.50 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

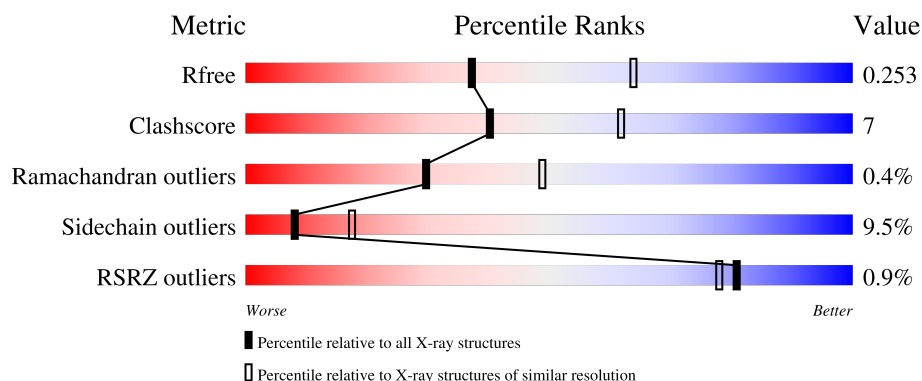
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

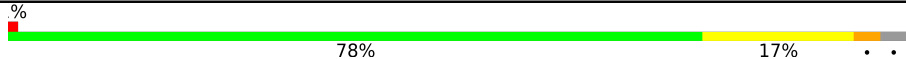
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	436	 78% 17% . .
1	B	436	 76% 17% . .

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
4	PEG	B	509	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6432 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 Serine Hydrolase OSH18.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	Se	0	0	0
			3161	1993	539	614	1	14			
1	B	426	Total	C	N	O	S	Se	0	0	0
			3178	2000	540	622	1	15			

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	11	Total	Na	0	0
			11	11		
2	B	7	Total	Na	0	0
			7	7		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			7	4	3		

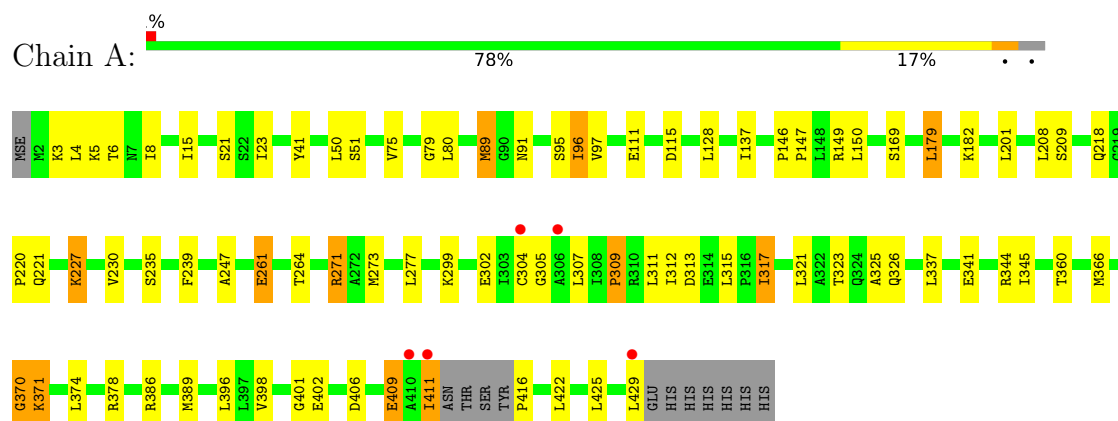
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	38	Total	O	0	1
			39	39		
5	B	28	Total	O	0	0
			28	28		

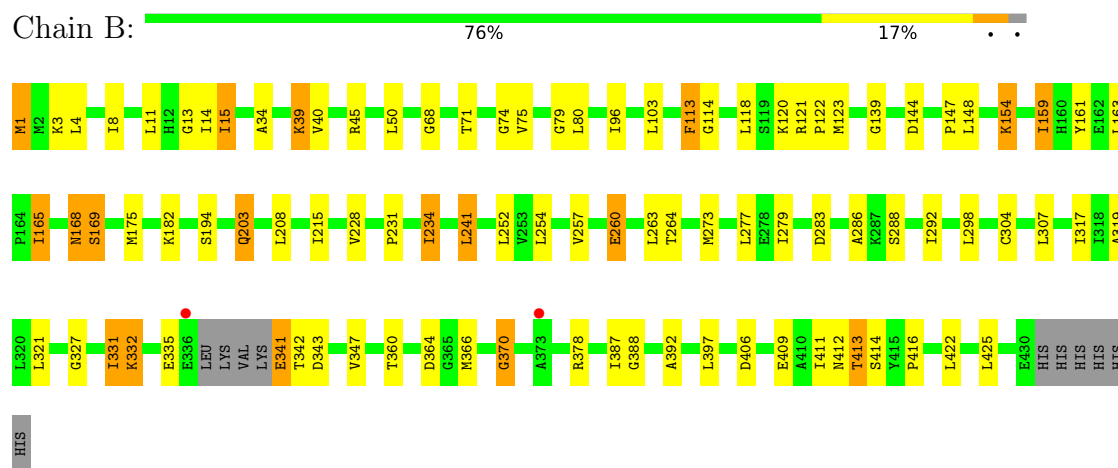
3 Residue-property plots

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: Serine Hydrolase OSH18



• Molecule 1: Serine Hydrolase OSH18



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	73.12Å 73.12Å 163.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.74 – 2.50 19.74 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (19.74-2.50) 99.0 (19.74-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.27 (at 2.50Å)	Xtriage
Refinement program	PHENIX dev_1269	Depositor
R, R_{free}	0.188 , 0.253 0.188 , 0.253	Depositor DCC
R_{free} test set	1478 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.052 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6432	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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: PEG, CL, NA

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	0.57	0/3185	0.90	7/4282 (0.2%)
1	B	0.53	0/3202	0.90	3/4306 (0.1%)
All	All	0.55	0/6387	0.90	10/8588 (0.1%)

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	B	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	370	GLY	N-CA-C	8.90	126.84	112.06
1	A	401	GLY	N-CA-C	-6.16	105.76	112.04
1	A	128	LEU	CA-C-N	-6.15	112.67	119.19
1	A	128	LEU	C-N-CA	-6.15	112.67	119.19
1	A	409	GLU	N-CA-C	-5.36	101.97	109.96
1	A	247	ALA	CA-C-N	5.14	125.04	119.85
1	A	247	ALA	C-N-CA	5.14	125.04	119.85
1	B	13	GLY	N-CA-C	5.10	117.49	110.56
1	B	283	ASP	CA-C-N	5.01	124.80	119.28
1	B	283	ASP	C-N-CA	5.01	124.80	119.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	169	SER	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	3161	0	3274	36	1
1	B	3178	0	3272	49	1
2	A	11	0	0	0	0
2	B	7	0	0	0	0
3	B	1	0	0	0	0
4	B	7	0	10	6	0
5	A	39	0	0	0	0
5	B	28	0	0	0	0
All	All	6432	0	6556	85	1

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 7.

All (85) 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:89:MSE:HG3	1:A:96:ILE:HG12	1.61	0.82
1:B:203:GLN:HE22	4:B:509:PEG:H11	1.43	0.81
1:B:327:GLY:H	1:B:370:GLY:HA2	1.46	0.80
1:B:175:MSE:HE2	1:B:215:ILE:HG21	1.64	0.79
1:B:241:LEU:HD13	1:B:254:LEU:HD12	1.69	0.73
1:A:239:PHE:HA	1:A:317:ILE:HG12	1.71	0.73
1:B:15:ILE:HD11	1:B:425:LEU:HD11	1.74	0.69
1:A:345:ILE:HG23	1:A:366:MSE:HE3	1.76	0.66
1:A:271:ARG:NH1	1:A:277:LEU:O	2.29	0.66
1:B:203:GLN:NE2	4:B:509:PEG:H11	2.11	0.65
1:B:159:ILE:HD13	1:B:161:TYR:HB2	1.78	0.65
1:A:304:CYS:HB2	1:A:307:LEU:HD12	1.79	0.65
1:A:273:MSE:HE3	1:A:325:ALA:HB2	1.80	0.64
1:A:182:LYS:HD2	1:A:220:PRO:HD3	1.80	0.62
1:B:319:ALA:HB2	1:B:366:MSE:HE1	1.82	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:273:MSE:HE3	1:B:298:LEU:HB3	1.83	0.60
1:B:252:LEU:HD22	1:B:425:LEU:HD22	1.84	0.60
1:B:304:CYS:HB3	1:B:307:LEU:HD12	1.83	0.59
1:A:305:GLY:HA2	1:A:311:LEU:HD22	1.84	0.59
1:B:228:VAL:HG13	4:B:509:PEG:H22	1.84	0.59
1:A:179:LEU:HD21	1:A:201:LEU:HD13	1.85	0.58
1:A:8:ILE:HD13	1:A:396:LEU:HD22	1.86	0.57
1:A:3:LYS:HD2	1:A:402:GLU:HB3	1.87	0.57
1:B:114:GLY:HA3	1:B:118:LEU:HD22	1.86	0.56
1:A:89:MSE:HG3	1:A:96:ILE:CG1	2.32	0.56
1:A:15:ILE:HD11	1:A:425:LEU:HD11	1.86	0.56
1:B:39:LYS:HD2	1:B:71:THR:HG23	1.88	0.56
1:B:286:ALA:HB2	4:B:509:PEG:H12	1.86	0.55
1:B:332:LYS:HG3	1:B:364:ASP:CG	2.32	0.55
1:B:120:LYS:O	1:B:122:PRO:HD3	2.07	0.54
1:A:137:ILE:HD12	1:A:150:LEU:HD12	1.87	0.54
1:A:218:GLN:O	1:A:221:GLN:NE2	2.33	0.53
1:A:325:ALA:O	1:A:370:GLY:HA3	2.08	0.53
1:A:91:ASN:HB3	1:A:115:ASP:OD1	2.08	0.53
1:B:121:ARG:HH22	1:B:168:ASN:HD22	1.58	0.51
1:A:6:THR:HG23	1:A:398:VAL:HB	1.93	0.51
1:A:41:TYR:HE1	1:A:227:LYS:HG2	1.77	0.50
1:A:273:MSE:HE1	1:A:321:LEU:HG	1.93	0.50
1:B:286:ALA:O	4:B:509:PEG:H41	2.11	0.50
1:A:111:GLU:HG2	1:A:149:ARG:HG3	1.94	0.50
1:A:75:VAL:O	1:A:79:GLY:HA3	2.12	0.50
1:B:332:LYS:NZ	1:B:364:ASP:OD2	2.30	0.49
1:A:239:PHE:CE1	1:A:389:MSE:HB3	2.48	0.49
1:A:312:ILE:HG21	1:A:341:GLU:HG2	1.94	0.48
1:B:343:ASP:O	1:B:347:VAL:HG23	2.13	0.48
1:A:227:LYS:HE2	1:A:227:LYS:HB3	1.72	0.48
1:B:113:PHE:HB3	1:B:147:PRO:HB3	1.96	0.47
1:B:388:GLY:C	1:B:411:ILE:HD11	2.40	0.47
1:B:286:ALA:CB	4:B:509:PEG:H12	2.44	0.47
1:A:411:ILE:HG13	1:A:416:PRO:HA	1.97	0.47
1:B:331:ILE:HD12	1:B:366:MSE:HE3	1.96	0.47
1:B:4:LEU:HD13	1:B:392:ALA:HA	1.97	0.46
1:A:396:LEU:HD11	1:A:422:LEU:HD11	1.98	0.46
1:B:327:GLY:O	1:B:370:GLY:N	2.49	0.45
1:B:341:GLU:N	1:B:342:THR:HA	2.31	0.45
1:B:378:ARG:HD2	1:B:406:ASP:OD2	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:409:GLU:O	1:B:413:THR:HG22	2.16	0.44
1:B:182:LYS:HE2	1:B:182:LYS:HB3	1.86	0.44
1:B:1:MSE:HE2	1:B:378:ARG:CZ	2.48	0.44
1:B:75:VAL:O	1:B:79:GLY:HA3	2.17	0.44
1:B:96:ILE:HG22	1:B:123:MSE:HE2	1.99	0.44
1:B:139:GLY:N	1:B:144:ASP:OD1	2.43	0.43
1:B:45:ARG:NH2	1:B:68:GLY:HA2	2.33	0.43
1:A:371:LYS:HA	1:A:371:LYS:HD3	1.74	0.43
1:B:234:ILE:CD1	1:B:260:GLU:HG3	2.48	0.43
1:B:165:ILE:HD13	1:B:165:ILE:HA	1.73	0.43
1:B:175:MSE:CE	1:B:215:ILE:HD13	2.48	0.43
1:B:277:LEU:HA	1:B:292:ILE:O	2.19	0.43
1:A:313:ASP:HA	1:A:344:ARG:NH1	2.34	0.43
1:A:299:LYS:HD3	1:A:326:GLN:HB3	2.01	0.42
1:A:305:GLY:CA	1:A:311:LEU:HD22	2.50	0.42
1:A:97:VAL:HG11	1:A:169:SER:HB2	2.02	0.41
1:B:273:MSE:HE1	1:B:321:LEU:HD11	2.03	0.41
1:B:154:LYS:H	1:B:154:LYS:HG3	1.67	0.41
1:B:414:SER:O	1:B:416:PRO:HD3	2.19	0.41
1:B:231:PRO:HB3	1:B:288:SER:HB3	2.03	0.41
1:B:387:ILE:HD13	1:B:387:ILE:HA	1.94	0.41
1:A:309:PRO:C	1:A:311:LEU:H	2.28	0.41
1:B:1:MSE:HE2	1:B:378:ARG:NH2	2.36	0.41
1:B:8:ILE:HD12	1:B:11:LEU:HD13	2.03	0.41
1:B:343:ASP:OD2	1:B:343:ASP:N	2.54	0.40
1:B:34:ALA:O	1:B:74:GLY:HA3	2.21	0.40
1:A:23:ILE:HG22	1:A:230:VAL:HG22	2.02	0.40
1:A:261:GLU:H	1:A:261:GLU:HG3	1.50	0.40
1:A:146:PRO:HA	1:A:147:PRO:C	2.46	0.40

All (1) 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)
1:A:271:ARG:NH2	1:B:279:ILE:O[1_565]	2.15	0.05

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	420/436 (96%)	401 (96%)	18 (4%)	1 (0%)	43	63
1	B	422/436 (97%)	402 (95%)	18 (4%)	2 (0%)	24	43
All	All	842/872 (97%)	803 (95%)	36 (4%)	3 (0%)	30	49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	370	GLY
1	B	168	ASN
1	A	309	PRO

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	346/343 (101%)	315 (91%)	31 (9%)	9	19
1	B	348/343 (102%)	313 (90%)	35 (10%)	7	15
All	All	694/686 (101%)	628 (90%)	66 (10%)	8	17

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	5	LYS
1	A	21	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	50	LEU
1	A	51	SER
1	A	80	LEU
1	A	89	MSE
1	A	95	SER
1	A	96	ILE
1	A	179	LEU
1	A	208	LEU
1	A	209	SER
1	A	227	LYS
1	A	235	SER
1	A	261	GLU
1	A	264	THR
1	A	271	ARG
1	A	302	GLU
1	A	315	LEU
1	A	317	ILE
1	A	323	THR
1	A	337	LEU
1	A	360	THR
1	A	371	LYS
1	A	374	LEU
1	A	378	ARG
1	A	386	ARG
1	A	406	ASP
1	A	409	GLU
1	A	411	ILE
1	A	429	LEU
1	B	1	MSE
1	B	3	LYS
1	B	14	ILE
1	B	15	ILE
1	B	39	LYS
1	B	40	VAL
1	B	50	LEU
1	B	80	LEU
1	B	103	LEU
1	B	113	PHE
1	B	148	LEU
1	B	154	LYS
1	B	159	ILE
1	B	163	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	165	ILE
1	B	169	SER
1	B	194	SER
1	B	203	GLN
1	B	208	LEU
1	B	234	ILE
1	B	241	LEU
1	B	257	VAL
1	B	260	GLU
1	B	263	LEU
1	B	264	THR
1	B	317	ILE
1	B	331	ILE
1	B	332	LYS
1	B	335	GLU
1	B	341	GLU
1	B	360	THR
1	B	397	LEU
1	B	412	ASN
1	B	413	THR
1	B	422	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	346	GLN
1	B	54	GLN
1	B	180	GLN
1	B	195	ASN
1	B	202	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 19 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	PEG	B	509	-	6,6,6	0.68	0	5,5,5	0.49	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
4	PEG	B	509	-	-	0/4/4/4	-

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.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	509	PEG	6	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	410/436 (94%)	-0.32	5 (1%) 76 73	27, 46, 82, 108	0
1	B	411/436 (94%)	-0.15	2 (0%) 87 85	29, 56, 87, 118	0
All	All	821/872 (94%)	-0.24	7 (0%) 81 78	27, 50, 85, 118	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	429	LEU	2.7
1	B	373	ALA	2.6
1	A	306	ALA	2.5
1	A	410	ALA	2.3
1	A	304	CYS	2.2
1	B	336	GLU	2.0
1	A	411	ILE	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	NA	A	511	1/1	0.81	0.15	64,64,64,64	0
2	NA	B	507	1/1	0.88	0.10	64,64,64,64	0
2	NA	B	505	1/1	0.89	0.10	66,66,66,66	0
2	NA	B	504	1/1	0.90	0.13	49,49,49,49	0
4	PEG	B	509	7/7	0.90	0.13	67,71,79,86	0
2	NA	A	508	1/1	0.91	0.07	53,53,53,53	0
2	NA	A	506	1/1	0.91	0.08	75,75,75,75	0
2	NA	A	507	1/1	0.91	0.08	66,66,66,66	0
2	NA	B	502	1/1	0.92	0.11	53,53,53,53	0
2	NA	A	509	1/1	0.93	0.07	63,63,63,63	0
2	NA	B	503	1/1	0.93	0.15	52,52,52,52	0
2	NA	B	508	1/1	0.94	0.05	57,57,57,57	0
2	NA	A	504	1/1	0.94	0.06	49,49,49,49	0
2	NA	A	505	1/1	0.95	0.10	34,34,34,34	0
2	NA	A	501	1/1	0.95	0.10	57,57,57,57	0
2	NA	B	506	1/1	0.95	0.06	64,64,64,64	0
2	NA	A	502	1/1	0.96	0.12	55,55,55,55	0
2	NA	A	510	1/1	0.97	0.07	61,61,61,61	0
2	NA	A	503	1/1	0.97	0.05	45,45,45,45	0
3	CL	B	501	1/1	0.98	0.07	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.