



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 11:59 AM UTC

PDB ID : 8P1H / pdb_00008p1h
Title : Crystal structure of the chimera of human 14-3-3 zeta and phosphorylated cytoplasmic loop fragment of the alpha7 acetylcholine receptor
Authors : Boyko, K.M.; Kapitonova, A.A.; Tugaeva, K.V.; Varfolomeeva, L.A.; Lyukmanova, E.N.; Sluchanko, N.N.
Deposited on : 2023-05-12
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

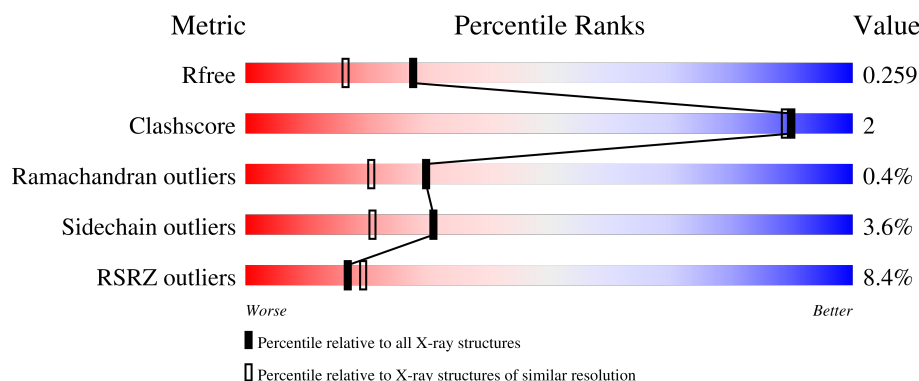
1 Overall quality at a glance

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)

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	248	10% 85% 11% ..
1	B	248	6% 91% 5% ..

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	AZI	A	302	-	X	-	-
3	AZI	A	303	-	X	-	-
3	AZI	A	304	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4214 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 Tyrosine 3-monooxygenase/tryptophan 5-monooxygenase activation protein zeta, Neuronal acetylcholine receptor subunit alpha-7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	P	S	0	7	0
			1907	1185	327	382	1	12			
1	B	240	Total	C	N	O	P	S	0	7	0
			1897	1181	328	377	1	10			

There are 44 discrepancies between the modelled and reference sequences:

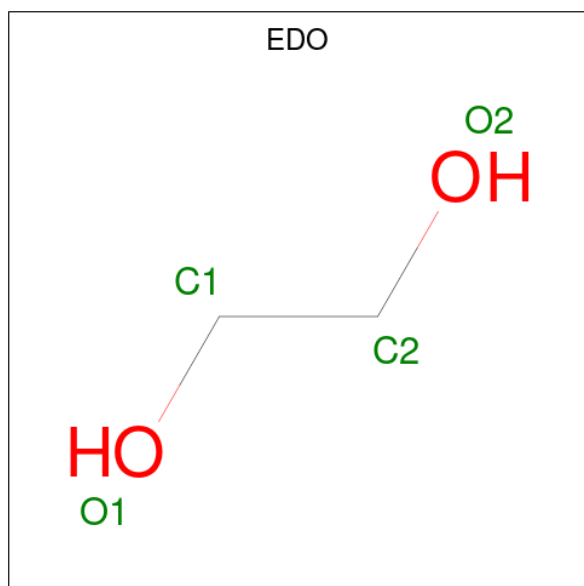
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP E7EX29
A	-1	PRO	-	expression tag	UNP E7EX29
A	0	HIS	-	expression tag	UNP E7EX29
A	58	ALA	SER	engineered mutation	UNP E7EX29
A	73	ALA	-	linker	UNP E7EX29
A	74	ALA	-	linker	UNP E7EX29
A	75	ALA	-	linker	UNP E7EX29
A	?	-	VAL	deletion	UNP E7EX29
A	?	-	ARG	deletion	UNP E7EX29
A	?	-	PHE	deletion	UNP E7EX29
A	?	-	GLY	deletion	UNP E7EX29
A	?	-	VAL	deletion	UNP E7EX29
A	?	-	SER	deletion	UNP E7EX29
A	?	-	TYR	deletion	UNP E7EX29
A	?	-	PRO	deletion	UNP E7EX29
A	157	ALA	-	linker	UNP E7EX29
A	158	ALA	-	linker	UNP E7EX29
A	159	ALA	-	linker	UNP E7EX29
A	230	GLY	-	linker	UNP E7EX29
A	231	GLY	-	linker	UNP E7EX29
A	232	GLY	-	linker	UNP E7EX29
A	233	GLY	-	linker	UNP E7EX29
B	-2	GLY	-	expression tag	UNP E7EX29
B	-1	PRO	-	expression tag	UNP E7EX29

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	HIS	-	expression tag	UNP E7EX29
B	58	ALA	SER	engineered mutation	UNP E7EX29
B	73	ALA	-	linker	UNP E7EX29
B	74	ALA	-	linker	UNP E7EX29
B	75	ALA	-	linker	UNP E7EX29
B	?	-	VAL	deletion	UNP E7EX29
B	?	-	ARG	deletion	UNP E7EX29
B	?	-	PHE	deletion	UNP E7EX29
B	?	-	GLY	deletion	UNP E7EX29
B	?	-	VAL	deletion	UNP E7EX29
B	?	-	SER	deletion	UNP E7EX29
B	?	-	TYR	deletion	UNP E7EX29
B	?	-	PRO	deletion	UNP E7EX29
B	157	ALA	-	linker	UNP E7EX29
B	158	ALA	-	linker	UNP E7EX29
B	159	ALA	-	linker	UNP E7EX29
B	230	GLY	-	linker	UNP E7EX29
B	231	GLY	-	linker	UNP E7EX29
B	232	GLY	-	linker	UNP E7EX29
B	233	GLY	-	linker	UNP E7EX29

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



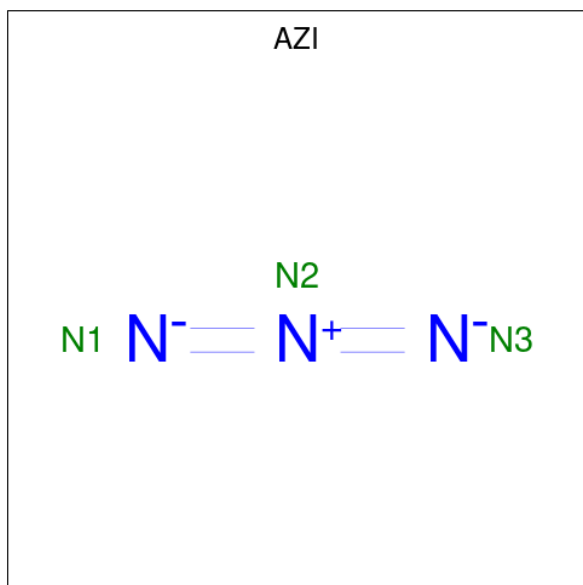
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

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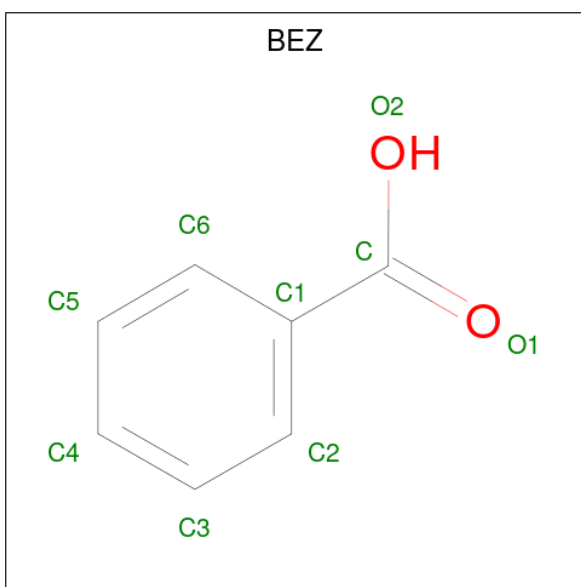
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is AZIDE ION (CCD ID: AZI) (formula: N₃).



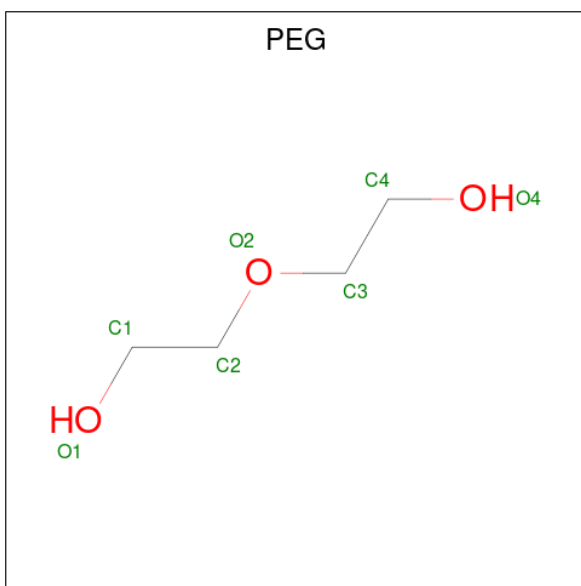
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	N	0	0
			3	3		
3	A	1	Total	N	0	0
			3	3		
3	A	1	Total	N	0	0
			3	3		

- Molecule 4 is BENZOIC ACID (CCD ID: BEZ) (formula: C₇H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			9	7	2		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

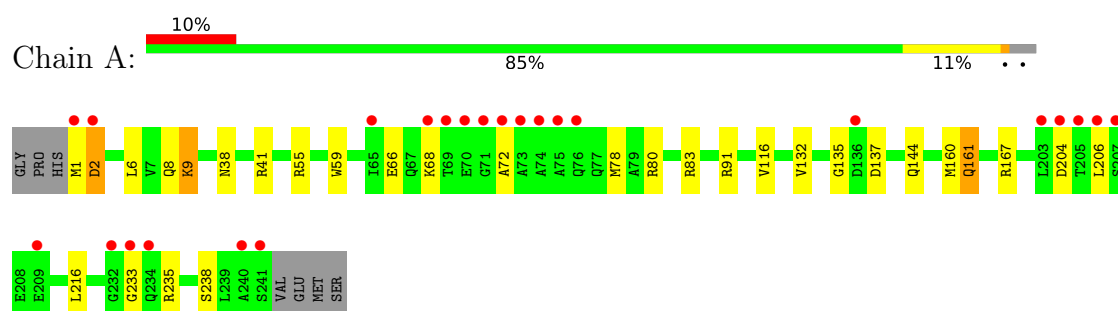
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	177	Total 177	O 177	0	0
6	B	192	Total 192	O 192	0	0

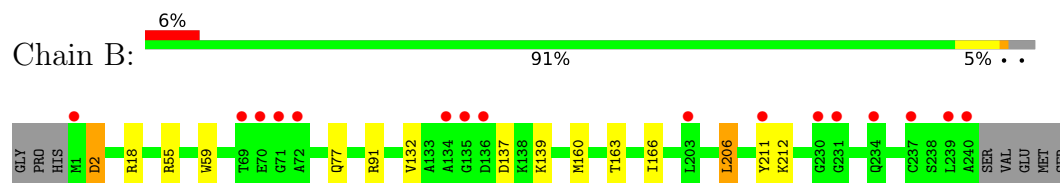
3 Residue-property plots [i](#)

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: Tyrosine 3-monooxygenase/tryptophan 5-monooxygenase activation protein zeta, Neuronal acetylcholine receptor subunit alpha-7



- Molecule 1: Tyrosine 3-monooxygenase/tryptophan 5-monooxygenase activation protein zeta, Neuronal acetylcholine receptor subunit alpha-7



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.63Å 103.12Å 112.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.06 – 1.95 21.06 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.7 (21.06-1.95) 97.7 (21.06-1.95)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0411	Depositor
R, R_{free}	0.209 , 0.253 0.216 , 0.259	Depositor DCC
R_{free} test set	2972 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4214	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.65% 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: SEP, AZI, BEZ, EDO, PEG

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	0.73	0/1946	1.10	1/2618 (0.0%)
1	B	0.72	0/1938	1.08	1/2609 (0.0%)
All	All	0.72	0/3884	1.09	2/5227 (0.0%)

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	7
1	B	0	2
All	All	0	9

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	2	ASP	CB-CA-C	5.80	119.42	109.51
1	A	2	ASP	CB-CA-C	5.61	119.11	109.51

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	135	GLY	Peptide
1	A	167	ARG	Sidechain
1	A	233	GLY	Peptide
1	A	235[A]	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	55	ARG	Sidechain
1	A	80	ARG	Sidechain
1	A	91	ARG	Sidechain
1	B	55	ARG	Sidechain
1	B	91	ARG	Sidechain

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	1907	0	1855	8	0
1	B	1897	0	1867	6	0
2	A	12	0	18	0	0
2	B	4	0	6	0	0
3	A	9	0	0	0	0
4	A	9	0	5	0	0
5	B	7	0	10	1	0
6	A	177	0	0	0	0
6	B	192	0	0	0	0
All	All	4214	0	3761	14	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 2.

All (14) 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:38[A]:ASN:OD1	1:A:41[A]:ARG:NH1	2.17	0.76
1:B:160:MET:HE1	1:B:166:ILE:HB	1.85	0.58
1:A:161:GLN:HE21	1:A:161:GLN:HA	1.73	0.52
1:A:116:VAL:HG11	1:A:160:MET:HE1	1.96	0.46
1:B:77[A]:GLN:OE1	1:B:77[A]:GLN:HA	2.14	0.46
1:B:163:THR:HB	1:B:206:LEU:CD1	2.48	0.44
1:B:206:LEU:HG	1:B:211:TYR:HA	1.99	0.44
1:B:59:TRP:CE2	1:B:132:VAL:HG12	2.55	0.42
1:A:66:GLU:OE2	1:A:83:ARG:NE	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ARG:HB3	5:B:302:PEG:H11	2.01	0.42
1:A:1:MET:HE3	1:A:6:LEU:HD21	2.02	0.41
1:A:59:TRP:CE2	1:A:132:VAL:HG12	2.55	0.41
1:A:1:MET:SD	1:A:9:LYS:HE3	2.61	0.41
1:A:78:MET:HE2	1:A:78:MET:HB3	1.97	0.41

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	245/248 (99%)	239 (98%)	4 (2%)	2 (1%)	16	8
1	B	245/248 (99%)	240 (98%)	5 (2%)	0	100	100
All	All	490/496 (99%)	479 (98%)	9 (2%)	2 (0%)	30	21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	ALA
1	A	68	LYS

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	200/204 (98%)	191 (96%)	9 (4%)	24	14
1	B	198/204 (97%)	193 (98%)	5 (2%)	42	34
All	All	398/408 (98%)	384 (96%)	14 (4%)	31	22

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	8	GLN
1	A	9	LYS
1	A	137	ASP
1	A	144	GLN
1	A	161	GLN
1	A	204	ASP
1	A	206	LEU
1	A	216	LEU
1	B	2	ASP
1	B	137	ASP
1	B	139	LYS
1	B	206	LEU
1	B	212	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	147	GLN
1	A	161	GLN
1	B	32	GLN
1	B	67	GLN
1	B	219	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	$\# Z > 2$
1	SEP	B	238	1	8,9,10	1.00	0	7,12,14	0.67	0
1	SEP	A	238	1	8,9,10	1.22	1 (12%)	7,12,14	1.19	1 (14%)

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
1	SEP	B	238	1	-	0/6/8/10	-
1	SEP	A	238	1	-	0/6/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	238	SEP	P-OG	-2.13	1.53	1.60

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	238	SEP	OG-CB-CA	2.03	110.12	108.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

9 ligands 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	EDO	A	301	-	3,3,3	0.36	0	2,2,2	1.53	0
3	AZI	A	304	-	2,2,2	3.59	2 (100%)	0,1,1	-	-
2	EDO	B	301	-	3,3,3	0.28	0	2,2,2	0.24	0
4	BEZ	A	307	-	9,9,9	1.00	1 (11%)	11,11,11	0.70	0
2	EDO	A	306	-	3,3,3	0.27	0	2,2,2	0.44	0
3	AZI	A	302	-	2,2,2	4.58	2 (100%)	0,1,1	-	-
3	AZI	A	303	-	2,2,2	3.91	2 (100%)	0,1,1	-	-
5	PEG	B	302	-	6,6,6	0.27	0	5,5,5	0.42	0
2	EDO	A	305	-	3,3,3	0.72	0	2,2,2	0.23	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
2	EDO	A	301	-	-	1/1/1/1	-
2	EDO	B	301	-	-	0/1/1/1	-
4	BEZ	A	307	-	-	0/4/4/4	0/1/1/1
2	EDO	A	306	-	-	0/1/1/1	-
5	PEG	B	302	-	-	2/4/4/4	-
2	EDO	A	305	-	-	1/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	AZI	N1-N2	-4.80	1.13	1.23
3	A	302	AZI	N3-N2	-4.34	1.14	1.23
3	A	303	AZI	N1-N2	-4.16	1.14	1.23
3	A	304	AZI	N1-N2	-3.65	1.15	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	303	AZI	N3-N2	-3.64	1.15	1.23
3	A	304	AZI	N3-N2	-3.53	1.15	1.23
4	A	307	BEZ	O2-C	-2.37	1.23	1.30

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	302	PEG	O1-C1-C2-O2
2	A	305	EDO	O1-C1-C2-O2
5	B	302	PEG	C4-C3-O2-C2
2	A	301	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	302	PEG	1	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	240/248 (96%)	0.30	24 (10%)	12 14	18, 32, 76, 114	8 (3%)
1	B	239/248 (96%)	0.30	16 (6%)	24 27	13, 33, 66, 97	8 (3%)
All	All	479/496 (96%)	0.30	40 (8%)	17 19	13, 33, 70, 114	16 (3%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	240	ALA	11.7
1	B	1	MET	5.6
1	A	234[A]	GLN	5.5
1	A	241	SER	5.4
1	A	75	ALA	5.2
1	A	73	ALA	4.8
1	A	1	MET	4.8
1	A	72	ALA	4.7
1	A	74	ALA	4.3
1	A	71	GLY	3.8
1	B	72	ALA	3.5
1	A	233	GLY	3.4
1	B	70	GLU	3.4
1	B	239	LEU	3.3
1	B	69	THR	3.2
1	B	71	GLY	3.1
1	A	69	THR	3.1
1	B	135	GLY	3.0
1	A	70	GLU	3.0
1	B	203	LEU	2.9
1	A	205	THR	2.8
1	B	234	GLN	2.7
1	B	136	ASP	2.6
1	A	207	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	240	ALA	2.5
1	A	206	LEU	2.4
1	A	68	LYS	2.4
1	A	209	GLU	2.4
1	B	211	TYR	2.4
1	B	237	CYS	2.4
1	A	203	LEU	2.4
1	A	2	ASP	2.3
1	A	65	ILE	2.3
1	A	76	GLN	2.3
1	B	230	GLY	2.3
1	B	231	GLY	2.3
1	A	232	GLY	2.2
1	B	134	ALA	2.2
1	A	204	ASP	2.2
1	A	136	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	SEP	A	238	10/11	0.98	0.05	25,27,33,34	0
1	SEP	B	238	10/11	0.98	0.05	27,29,39,45	0

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
3	AZI	A	304	3/3	0.84	0.15	46,46,59,61	0
5	PEG	B	302	7/7	0.87	0.20	40,53,67,75	0
2	EDO	A	305	4/4	0.88	0.24	50,51,53,56	0
2	EDO	B	301	4/4	0.90	0.12	47,51,54,56	0
3	AZI	A	303	3/3	0.92	0.09	36,36,48,51	0
2	EDO	A	301	4/4	0.92	0.12	36,37,39,41	0
3	AZI	A	302	3/3	0.92	0.09	42,42,51,51	0
2	EDO	A	306	4/4	0.93	0.13	47,47,49,53	0
4	BEZ	A	307	9/9	0.98	0.04	24,26,27,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.