

Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 12:18 PM EST

PDB ID : 5EFN

Title : Crystal structure of Danio rerio histone deacetylase 6 catalytic domain 2

(H574A) in complex with histone H4 Lys6 tripeptide substrate

Authors: Hai, Y.; Christianson, D.W.

Deposited on : 2015-10-24

Resolution : 1.80 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 (i) 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 (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

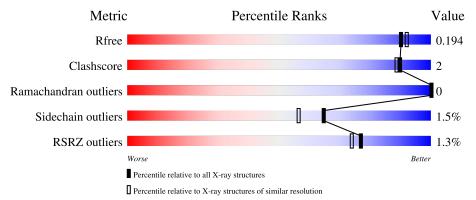
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.80 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	364	93%					
1	В	364	95%					
2	Е	3	67% 33%					
2	F	3	67% 67% 33%					



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 6396 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 Hdac6 protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	В	357	Total 2844	C 1796	N 502	O 527	S 19	0	13	0
1	A	357	Total 2830	C 1781	N 504	O 527	S 18	0	11	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	435	SER	-	expression tag	UNP A7YT55
В	436	ASN	-	expression tag	UNP A7YT55
В	437	ALA	-	expression tag	UNP A7YT55
В	438	GLY	-	expression tag	UNP A7YT55
В	439	GLY	-	expression tag	UNP A7YT55
В	574	ALA	HIS	engineered mutation	UNP A7YT55
A	435	SER	-	expression tag	UNP A7YT55
A	436	ASN	-	expression tag	UNP A7YT55
A	437	ALA	-	expression tag	UNP A7YT55
A	438	GLY	-	expression tag	UNP A7YT55
A	439	GLY	-	expression tag	UNP A7YT55
A	574	ALA	HIS	engineered mutation	UNP A7YT55

• Molecule 2 is a protein called histone H4 tripeptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	3	Total 22	C 13		O 5	0	0	0
2	Е	3	Total 22	C 13		O 5	0	0	0

• Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

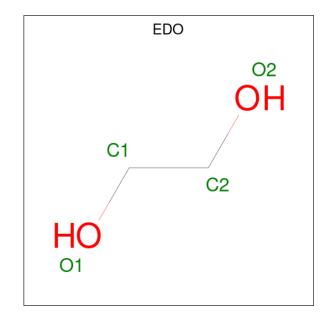


Mo	ol	Chain	Residues	Atoms	ZeroOcc	AltConf
3		В	1	Total Zn 1 1	0	0
3		A	1	Total Zn 1 1	0	0

• Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	2	Total K 2 2	0	0
4	A	2	Total K 2 2	0	0

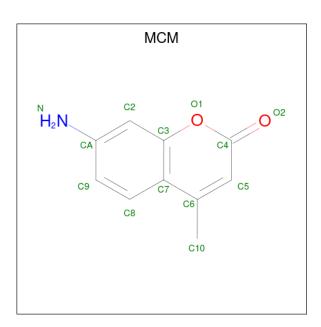
• Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

• Molecule 6 is 7-AMINO-4-METHYL-CHROMEN-2-ONE (three-letter code: MCM) (formula: $C_{10}H_9NO_2$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	F	1	Total 13				0	0
6	Е	1	Total 13	_	N 1	_	0	0

• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	В	344	Total O 344 344	0	0
7	A	288	Total O 288 288	0	0
7	Е	2	Total O 2 2	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: Hdac6 protein

Chain B:

95%

• Molecule 1: Hdac6 protein

Chain A:

93%

• Molecule 2: histone H4 tripeptide

67%

Chain F:

67%

67%

Chain E:

67%

Chain E:

67%

33%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	55.00Å 83.91Å 86.90Å	Donositor
a, b, c, α , β , γ	90.00° 98.12° 90.00°	Depositor
Resolution (Å)	12.34 - 1.80	Depositor
rtesolution (A)	12.34 - 1.80	EDS
% Data completeness	98.9 (12.34-1.80)	Depositor
(in resolution range)	98.9 (12.34-1.80)	EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.38 (at 1.80Å)	Xtriage
Refinement program	PHENIX 1.8.3_1479	Depositor
D D.	0.162 , 0.194	Depositor
R, R_{free}	0.164 , 0.194	DCC
R_{free} test set	3591 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.45, 59.0	EDS
L-test for twinning ²	$ < L > = 0.48, < L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6396	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 15.29% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: ZN, MCM, K, 5OL, EDO

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		
IVIOI	Chain	RMSZ # Z > 5		RMSZ	# Z > 5	
1	A	0.33	0/2924	0.48	0/3964	
1	В	0.35	0/2955	0.50	0/4007	
2	Е	0.14	0/8	0.27	0/9	
2	F	0.17	0/8	0.30	0/9	
All	All	0.34	0/5895	0.49	0/7989	

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	2830	0	2765	11	0
1	В	2844	0	2790	6	0
2	Е	22	0	4	1	0
2	F	22	0	4	1	0
3	A	1	0	0	0	0
3	В	1	0	0	0	0
4	A	2	0	0	0	0
4	В	2	0	0	0	0
5	A	4	0	6	0	0

Continued on next page...



Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	В	8	0	12	0	0
6	Е	13	0	8	1	0
6	F	13	0	8	2	0
7	A	288	0	0	5	0
7	В	344	0	0	2	0
7	Е	2	0	0	0	0
All	All	6396	0	5597	20	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 (20) 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:788:ARG:NH1	7:A:905:HOH:O	2.37	0.57
1:A:708:ARG:NH1	7:A:901:HOH:O	2.28	0.57
1:A:520:ARG:NH1	1:A:521:ASP:OD1	2.43	0.52
1:A:684:HIS:NE2	7:A:904:HOH:O	2.35	0.50
1:B:671:GLY:N	7:B:903:HOH:O	2.36	0.50
1:A:520:ARG:NH2	7:A:909:HOH:O	2.44	0.48
1:A:443:ILE:HG23	1:A:561:VAL:HG12	1.95	0.47
2:E:4:5OL:O	6:E:101:MCM:H2	2.15	0.47
1:A:672[B]:LYS:HD2	1:A:714:GLY:O	2.15	0.47
1:B:459[A]:TRP:CZ2	1:B:523:ASN:HB2	2.50	0.47
1:A:682:PHE:HA	1:A:686:VAL:HB	1.98	0.45
2:F:4:5OL:O	6:F:101:MCM:H2	2.18	0.43
1:B:565[A]:VAL:HG13	1:B:761:LEU:HD12	2.00	0.43
1:B:518:LYS:NZ	7:B:915:HOH:O	2.52	0.42
1:B:518:LYS:HA	1:B:518:LYS:HD3	1.94	0.42
1:A:443:ILE:HG13	7:A:1032:HOH:O	2.19	0.42
1:A:673:MET:SD	1:A:677:GLU:HG2	2.60	0.42
1:A:746:ASN:HB3	1:A:749[A]:SER:HB2	2.04	0.40
1:B:515[B]:GLU:HG3	1:B:516:HIS:CD2	2.57	0.40
6:F:101:MCM:H8	6:F:101:MCM:H101	1.97	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	Perce	ntiles
1	A	366/364 (100%)	358 (98%)	8 (2%)	0	100	100
1	В	$368/364 \ (101\%)$	359 (98%)	9 (2%)	0	100	100
2	E	1/3 (33%)	1 (100%)	0	0	100	100
2	F	1/3 (33%)	1 (100%)	0	0	100	100
All	All	736/734 (100%)	719 (98%)	17 (2%)	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	308/304 (101%)	302 (98%)	6 (2%)	57 46
1	В	312/304 (103%)	308 (99%)	4 (1%)	69 62
All	All	620/608 (102%)	610 (98%)	10 (2%)	65 54

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	465	GLU
1	В	520	ARG
1	В	573	HIS
1	В	611	TRP
1	A	443	ILE

Continued on next page...



Continued from previous page...

Mol	Chain	Res	Type
1	A	465	GLU
1	A	573	HIS
1	A	611	TRP
1	A	672[A]	LYS
1	A	672[B]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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		Dag	Link	Bo	ond leng	ths	Bond angles			
MIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	5OL	F	4	3,2,6	10,12,13	1.78	3 (30%)	5,15,17	0.78	0
2	5OL	E	4	3,2,6	10,12,13	1.83	3 (30%)	5,15,17	1.48	1 (20%)

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	5OL	F	4	3,2,6	-	0/7/11/13	-
2	5OL	E	4	3,2,6	-	0/7/11/13	-

All (6) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(\text{\AA})$
2	Е	4	5OL	O01-C08	3.44	1.43	1.38
2	F	4	5OL	O01-C08	3.08	1.42	1.38
2	F	4	5OL	C06-N07	-2.78	1.43	1.46
2	Е	4	5OL	C06-N07	-2.76	1.43	1.46
2	F	4	5OL	O02-C08	2.56	1.41	1.38
2	Е	4	5OL	O02-C08	2.54	1.41	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
2	E	4	5OL	C06-N07-C08	-3.16	112.86	116.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	4	5OL	1	0
2	Е	4	5OL	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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	Trunc	Chain	Res	s Link	Bo	Bond lengths			Bond angles		
	Type	Chain	rtes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
5	EDO	A	804	-	3,3,3	0.43	0	2,2,2	0.43	0	
6	MCM	Е	101	2	12,14,14	1.87	2 (16%)	16,20,20	1.71	4 (25%)	



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
WIOI	Type				Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
5	EDO	В	804	-	3,3,3	0.45	0	2,2,2	0.23	0
6	MCM	F	101	2	12,14,14	1.90	2 (16%)	16,20,20	1.66	4 (25%)
5	EDO	В	805	-	3,3,3	0.44	0	2,2,2	0.37	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
5	EDO	A	804	-	-	1/1/1/1	-
6	MCM	Е	101	2	-	-	0/2/2/2
5	EDO	В	804	-	-	1/1/1/1	-
6	MCM	F	101	2	-	-	0/2/2/2
5	EDO	В	805	-	-	0/1/1/1	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
6	F	101	MCM	C7-C3	5.12	1.48	1.41
6	Е	101	MCM	C7-C3	4.94	1.48	1.41
6	Е	101	MCM	C6-C7	3.05	1.49	1.42
6	F	101	MCM	C6-C7	3.05	1.49	1.42

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}(^{o})$
6	Е	101	MCM	O1-C3-C2	3.95	120.71	116.03
6	F	101	MCM	O1-C3-C2	3.81	120.54	116.03
6	Е	101	MCM	C2-C3-C7	-2.88	119.85	123.05
6	F	101	MCM	C2-C3-C7	-2.85	119.89	123.05
6	Е	101	MCM	C9-CA-N	-2.33	116.57	120.91
6	Е	101	MCM	C8-C7-C3	2.27	119.13	116.50
6	F	101	MCM	C8-C7-C3	2.21	119.06	116.50
6	F	101	MCM	O1-C4-C5	2.05	121.88	119.27

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	В	804	EDO	O1-C1-C2-O2

Continued on next page...



Continued from previous page...

M	[ol	Chain	Res	Type	Atoms
,	5	A	804	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

\mathbf{Mol}	Chain	Res	Type	Clashes	Symm-Clashes
6	Ε	101	MCM	1	0
6	F	101	MCM	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 (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, 95^{th} 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>	$\# \mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	357/364 (98%)	-0.44	4 (1%) 80 78	7, 14, 29, 43	0
1	В	357/364 (98%)	-0.62	2 (0%) 89 87	6, 11, 22, 38	0
2	E	2/3~(66%)	2.75	1 (50%) 0 0	65, 65, 65, 76	0
2	F	2/3 (66%)	2.46	2 (100%) 0 0	64, 64, 64, 74	0
All	All	718/734 (97%)	-0.51	9 (1%) 77 74	6, 12, 26, 76	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	462	HIS	4.7
1	A	442	PRO	4.6
1	В	442	PRO	4.4
2	Е	2	ARG	3.6
1	A	771	HIS	3.4
2	F	2	ARG	2.5
2	F	3	GLY	2.4
1	В	520	ARG	2.1
1	A	672[A]	LYS	2.0

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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	5OL	F	4	13/14	0.87	0.15	16,23,28,28	0
2	5OL	Е	4	13/14	0.87	0.14	18,20,23,24	0



6.3 Carbohydrates (i)

There are no monosaccharides 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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
6	MCM	Е	101	13/13	0.82	0.17	25,31,36,38	0
5	EDO	В	804	4/4	0.83	0.29	16,18,19,20	4
6	MCM	F	101	13/13	0.84	0.20	30,35,40,42	0
5	EDO	В	805	4/4	0.85	0.25	34,35,35,35	4
5	EDO	A	804	4/4	0.89	0.15	31,31,32,32	4
4	K	A	803	1/1	0.99	0.04	12,12,12,12	0
3	ZN	A	801	1/1	1.00	0.04	18,18,18,18	1
4	K	В	802	1/1	1.00	0.02	9,9,9,9	0
4	K	В	803	1/1	1.00	0.02	9,9,9,9	0
4	K	A	802	1/1	1.00	0.03	10,10,10,10	0
3	ZN	В	801	1/1	1.00	0.04	16,16,16,16	1

6.5 Other polymers (i)

There are no such residues in this entry.

