wwPDB X-ray Structure Validation Summary Report (i)

May 22, 2020 – 04:36 pm BST

PDB ID	:	5LNL
Title	:	Crystal structure of Hsf 1608-1749 putative domain 1
Authors	:	Thomsen, M.; Wright, J.; Ridley, J.; Goldman, A.
Deposited on	:	2016-08-05
Resolution	:	3.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see references (1)) were used in the production of this report:

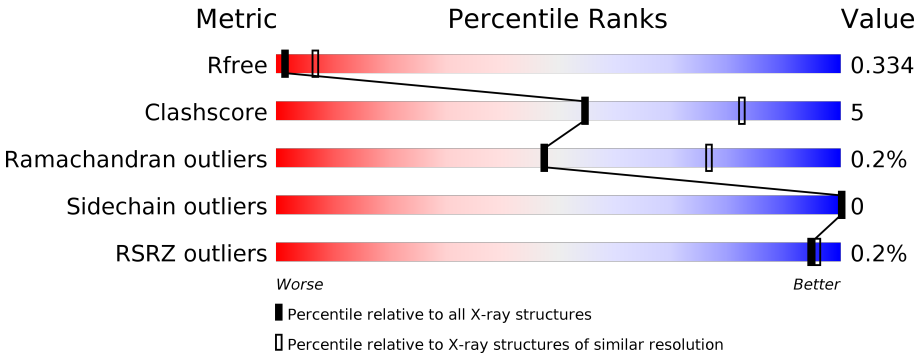
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 3.30 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}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 on 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	151	77% 9% 13%
1	B	151	72% 16% 13%
1	C	151	79% 9% 12%
1	D	151	74% 12% 15%
1	E	151	77% 10% 13%
1	F	151	% 74% 11% 15%

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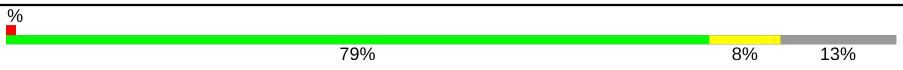
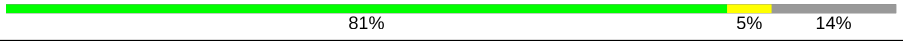
Mol	Chain	Length	Quality of chain
1	G	151	
1	H	151	
1	I	151	

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 8073 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 Hsf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			899	551	154	193	1			
1	B	132	Total	C	N	O		0	0	0
			907	556	157	194				
1	C	133	Total	C	N	O		0	0	0
			905	553	157	195				
1	D	129	Total	C	N	O		0	0	1
			883	541	153	189				
1	E	132	Total	C	N	O		0	0	0
			894	548	156	190				
1	F	129	Total	C	N	O		0	0	0
			893	546	152	195				
1	G	132	Total	C	N	O		0	0	0
			911	557	157	197				
1	H	130	Total	C	N	O		0	0	1
			883	542	154	187				
1	I	130	Total	C	N	O		0	0	0
			898	548	154	196				

There are 81 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1607	MET	-	initiating methionine	UNP P71401
A	1750	LEU	-	expression tag	UNP P71401
A	1751	GLU	-	expression tag	UNP P71401
A	1752	HIS	-	expression tag	UNP P71401
A	1753	HIS	-	expression tag	UNP P71401
A	1754	HIS	-	expression tag	UNP P71401
A	1755	HIS	-	expression tag	UNP P71401
A	1756	HIS	-	expression tag	UNP P71401
A	1757	HIS	-	expression tag	UNP P71401
B	1607	MET	-	initiating methionine	UNP P71401
B	1750	LEU	-	expression tag	UNP P71401

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1751	GLU	-	expression tag	UNP P71401
B	1752	HIS	-	expression tag	UNP P71401
B	1753	HIS	-	expression tag	UNP P71401
B	1754	HIS	-	expression tag	UNP P71401
B	1755	HIS	-	expression tag	UNP P71401
B	1756	HIS	-	expression tag	UNP P71401
B	1757	HIS	-	expression tag	UNP P71401
C	1607	MET	-	initiating methionine	UNP P71401
C	1750	LEU	-	expression tag	UNP P71401
C	1751	GLU	-	expression tag	UNP P71401
C	1752	HIS	-	expression tag	UNP P71401
C	1753	HIS	-	expression tag	UNP P71401
C	1754	HIS	-	expression tag	UNP P71401
C	1755	HIS	-	expression tag	UNP P71401
C	1756	HIS	-	expression tag	UNP P71401
C	1757	HIS	-	expression tag	UNP P71401
D	1607	MET	-	initiating methionine	UNP P71401
D	1750	LEU	-	expression tag	UNP P71401
D	1751	GLU	-	expression tag	UNP P71401
D	1752	HIS	-	expression tag	UNP P71401
D	1753	HIS	-	expression tag	UNP P71401
D	1754	HIS	-	expression tag	UNP P71401
D	1755	HIS	-	expression tag	UNP P71401
D	1756	HIS	-	expression tag	UNP P71401
D	1757	HIS	-	expression tag	UNP P71401
E	1607	MET	-	initiating methionine	UNP P71401
E	1750	LEU	-	expression tag	UNP P71401
E	1751	GLU	-	expression tag	UNP P71401
E	1752	HIS	-	expression tag	UNP P71401
E	1753	HIS	-	expression tag	UNP P71401
E	1754	HIS	-	expression tag	UNP P71401
E	1755	HIS	-	expression tag	UNP P71401
E	1756	HIS	-	expression tag	UNP P71401
E	1757	HIS	-	expression tag	UNP P71401
F	1607	MET	-	initiating methionine	UNP P71401
F	1750	LEU	-	expression tag	UNP P71401
F	1751	GLU	-	expression tag	UNP P71401
F	1752	HIS	-	expression tag	UNP P71401
F	1753	HIS	-	expression tag	UNP P71401
F	1754	HIS	-	expression tag	UNP P71401
F	1755	HIS	-	expression tag	UNP P71401
F	1756	HIS	-	expression tag	UNP P71401

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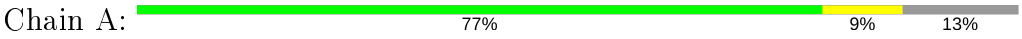
Chain	Residue	Modelled	Actual	Comment	Reference
F	1757	HIS	-	expression tag	UNP P71401
G	1607	MET	-	initiating methionine	UNP P71401
G	1750	LEU	-	expression tag	UNP P71401
G	1751	GLU	-	expression tag	UNP P71401
G	1752	HIS	-	expression tag	UNP P71401
G	1753	HIS	-	expression tag	UNP P71401
G	1754	HIS	-	expression tag	UNP P71401
G	1755	HIS	-	expression tag	UNP P71401
G	1756	HIS	-	expression tag	UNP P71401
G	1757	HIS	-	expression tag	UNP P71401
H	1607	MET	-	initiating methionine	UNP P71401
H	1750	LEU	-	expression tag	UNP P71401
H	1751	GLU	-	expression tag	UNP P71401
H	1752	HIS	-	expression tag	UNP P71401
H	1753	HIS	-	expression tag	UNP P71401
H	1754	HIS	-	expression tag	UNP P71401
H	1755	HIS	-	expression tag	UNP P71401
H	1756	HIS	-	expression tag	UNP P71401
H	1757	HIS	-	expression tag	UNP P71401
I	1607	MET	-	initiating methionine	UNP P71401
I	1750	LEU	-	expression tag	UNP P71401
I	1751	GLU	-	expression tag	UNP P71401
I	1752	HIS	-	expression tag	UNP P71401
I	1753	HIS	-	expression tag	UNP P71401
I	1754	HIS	-	expression tag	UNP P71401
I	1755	HIS	-	expression tag	UNP P71401
I	1756	HIS	-	expression tag	UNP P71401
I	1757	HIS	-	expression tag	UNP P71401

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

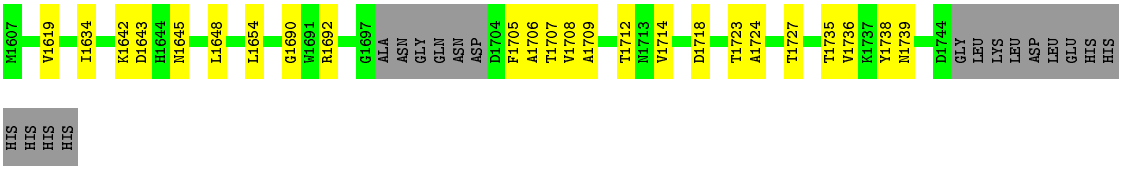
- Molecule 1: Hsf

Chain A:



- Molecule 1: Hsf

Chain B:

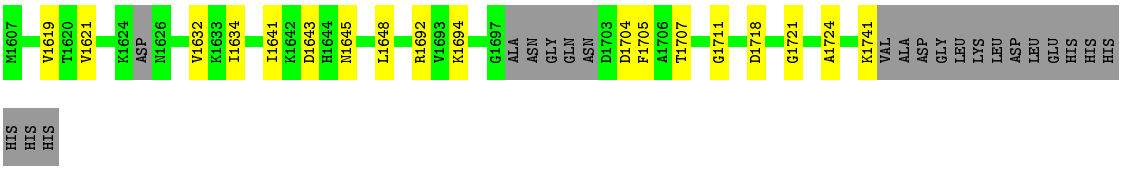


- Molecule 1: Hsf

Chain C:

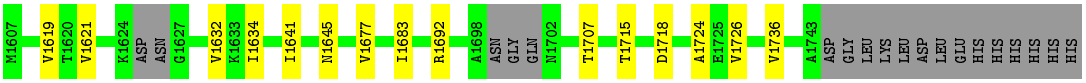
- Molecule 1: Hsf

Chain D:



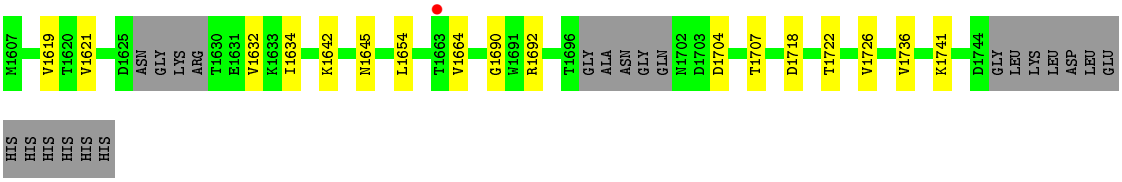
- Molecule 1: Hsf

Chain E:



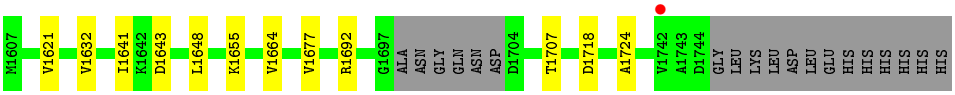
• Molecule 1: Hsf

Chain F:



• Molecule 1: Hsf

Chain G:

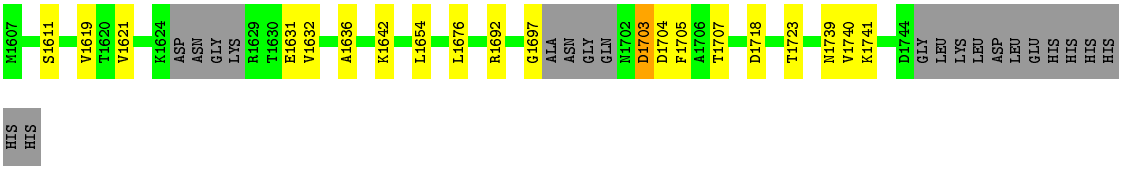


• Molecule 1: Hsf

Chain H:

• Molecule 1: Hsf

Chain I:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	128.44Å 50.40Å 256.80Å 90.00° 101.94° 90.00°	Depositor
Resolution (Å)	29.60 – 3.30 29.60 – 3.30	Depositor EDS
% Data completeness (in resolution range)	97.6 (29.60-3.30) 97.6 (29.60-3.30)	Depositor EDS
R_{merge}	0.33	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.44 (at 3.31Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.296 , 0.334 0.302 , 0.334	Depositor DCC
R_{free} test set	1179 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å^2)	56.8	Xtriage
Anisotropy	0.979	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.19 , -32.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.40$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	0.085 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.85	EDS
Total number of atoms	8073	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

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	0.21	0/904	0.41	0/1229
1	B	0.21	0/913	0.40	0/1243
1	C	0.21	0/911	0.41	0/1242
1	D	0.21	0/888	0.40	0/1206
1	E	0.21	0/899	0.39	0/1224
1	F	0.21	0/898	0.40	0/1224
1	G	0.21	0/917	0.40	0/1249
1	H	0.21	0/888	0.39	0/1209
1	I	0.21	0/903	0.39	0/1231
All	All	0.21	0/8121	0.40	0/11057

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

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	899	0	838	9	0
1	B	907	0	846	18	0
1	C	905	0	828	9	0
1	D	883	0	819	11	0
1	E	894	0	824	9	0
1	F	893	0	821	11	0
1	G	911	0	845	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	883	0	814	5	0
1	I	898	0	826	12	0
All	All	8073	0	7461	81	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 5.

The worst 5 of 81 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:1645:ASN:HB3	1:B:1692:ARG:HD2	1.74	0.69
1:A:1723:THR:HB	1:A:1739:ASN:HB2	1.76	0.68
1:A:1692:ARG:NH2	1:A:1707:THR:OG1	2.29	0.66
1:D:1692:ARG:NH1	1:D:1707:THR:OG1	2.29	0.65
1:B:1723:THR:HB	1:B:1739:ASN:HB2	1.79	0.63

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	125/151 (83%)	122 (98%)	3 (2%)	0	100	100
1	B	128/151 (85%)	125 (98%)	3 (2%)	0	100	100
1	C	129/151 (85%)	125 (97%)	3 (2%)	1 (1%)	19	51
1	D	123/151 (82%)	121 (98%)	2 (2%)	0	100	100
1	E	126/151 (83%)	124 (98%)	2 (2%)	0	100	100
1	F	123/151 (82%)	118 (96%)	5 (4%)	0	100	100
1	G	128/151 (85%)	126 (98%)	2 (2%)	0	100	100
1	H	124/151 (82%)	122 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	I	124/151 (82%)	122 (98%)	1 (1%)	1 (1%)	19 51
All	All	1130/1359 (83%)	1105 (98%)	23 (2%)	2 (0%)	47 77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	1703	ASP
1	C	1624	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	89/121 (74%)	89 (100%)	0	100 100
1	B	89/121 (74%)	89 (100%)	0	100 100
1	C	87/121 (72%)	87 (100%)	0	100 100
1	D	86/121 (71%)	86 (100%)	0	100 100
1	E	85/121 (70%)	85 (100%)	0	100 100
1	F	89/121 (74%)	89 (100%)	0	100 100
1	G	90/121 (74%)	90 (100%)	0	100 100
1	H	84/121 (69%)	84 (100%)	0	100 100
1	I	90/121 (74%)	90 (100%)	0	100 100
All	All	789/1089 (72%)	789 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Some 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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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

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, 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>2	OWAB(Å^2)	Q<0.9
1	A	131/151 (86%)	-0.53	0 100 100	29, 55, 95, 107	0
1	B	132/151 (87%)	-0.61	0 100 100	27, 51, 84, 101	0
1	C	133/151 (88%)	-0.60	0 100 100	30, 58, 88, 112	0
1	D	129/151 (85%)	-0.62	0 100 100	32, 58, 79, 105	0
1	E	132/151 (87%)	-0.55	0 100 100	35, 60, 94, 116	0
1	F	129/151 (85%)	-0.61	1 (0%) 86 86	30, 54, 100, 123	0
1	G	132/151 (87%)	-0.56	1 (0%) 86 86	25, 58, 87, 105	0
1	H	130/151 (86%)	-0.64	0 100 100	34, 59, 90, 99	0
1	I	130/151 (86%)	-0.59	0 100 100	34, 57, 83, 91	0
All	All	1178/1359 (86%)	-0.59	2 (0%) 95 96	25, 57, 90, 123	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	1742	VAL	5.0
1	F	1663	THR	2.4

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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.