

Electronic Supporting Information (ESI) for

**Eu<sub>2</sub>Pt<sub>3</sub>Pb<sub>5</sub> and SrPt<sub>2</sub>Pb<sub>4</sub> — lead-based intermetallics**

**with Y<sub>2</sub>Rh<sub>3</sub>Sn<sub>5</sub> / NdRh<sub>2</sub>Sn<sub>4</sub>-type polyanionic**

**platinum-lead 3D frameworks**

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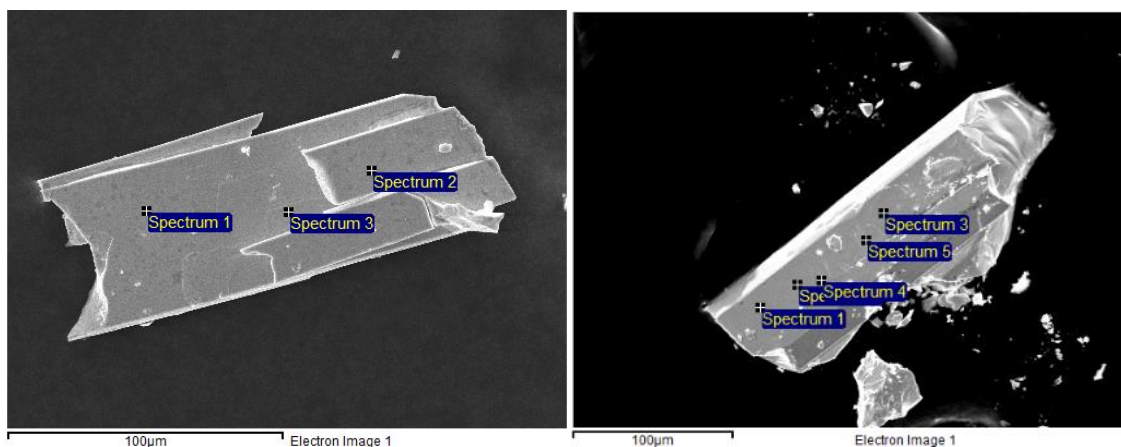
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**Figure S1.** SEM image of the  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$  (left) and  $\text{SrPt}_2\text{Pb}_4$  (right) crystals with analysis points marked.

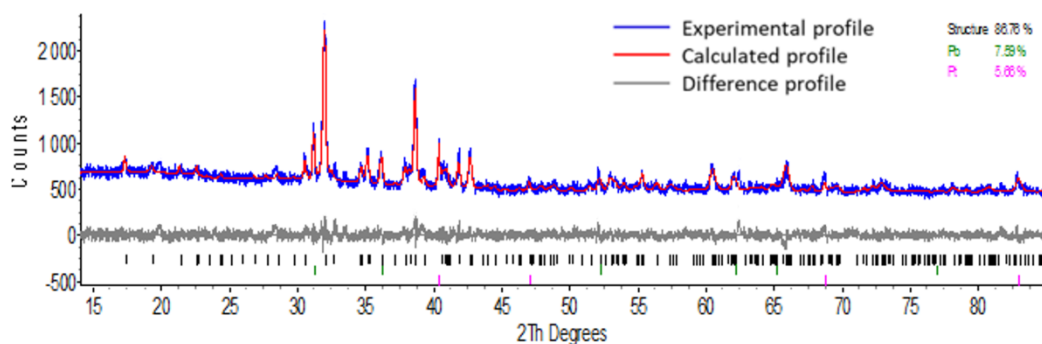
**Table S1.** Averaged results of the EDX analysis of the  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$  and  $\text{SrPt}_2\text{Pb}_4$  crystals.

System	Element	measurement by <i>ED</i> -spectrometer (atomic %)				Average formula
		Mean	Std. deviation	Max.	Min.	
Eu-Pt-Pb	Eu	20.03	0.92	20.87	18.70	$\text{Eu}_{2.00(9)}\text{Pt}_{2.99(5)}\text{P}_{5.00(7)}$
	Pt	29.93	0.50	30.43	29.13	
	Pb	50.05	0.74	50.97	48.97	
Sr-Pt-Pb	Sr	14.99	0.31	15.24	14.23	$\text{Sr}_{1.00(2)}\text{Pt}_{2.01(3)}\text{P}_{3.65(3)}$
	Pt	30.25	0.48	31.22	29.63	
	Pb	54.76	0.44	54.05	55.49	

**Table S2.** Anisotropic displacement parameters for  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$  and  $\text{SrPt}_2\text{Pb}_4$ .

$\text{Eu}_2\text{Pt}_3\text{Pb}_5$						
<i>Atom</i>	$U_{11}$	$U_{22}$	$U_{33}$	$U_{12}$	$U_{13}$	$U_{23}$
Pt1	0.0095(10)	0.0060(9)	0.0133(11)	0.0005(7)	0.0000	0.0000
Pt2	0.0081(8)	0.0042(8)	0.0157(10)	-0.0012(8)	0.0000	0.0000
Pt3	0.0093(10)	0.0051(8)	0.0148(10)	-0.0004(7)	0.0000	0.0000
Eu1	0.0102(10)	0.0055(10)	0.0147(12)	-0.0004(10)	0.0000	0.0000
Eu2	0.0085(11)	0.0051(10)	0.0144(14)	-0.0030(9)	0.0000	0.0000
Pb1	0.0082(10)	0.0036(9)	0.0149(9)	-0.0005(7)	0.0000	0.0000
Pb2	0.0122(10)	0.0077(9)	0.0161(9)	-0.0001(7)	0.0000	0.0000
Pb3	0.0086(9)	0.0038(8)	0.0157(10)	0.0005(6)	0.0000	0.0000
Pb4	0.0118(9)	0.0033(7)	0.0143(11)	-0.0007(7)	0.0000	0.0000
Pb5	0.0088(9)	0.0051(8)	0.0128(9)	0.0004(6)	0.0000	0.0000
$\text{SrPt}_2\text{Pb}_4$						

<i>Atom</i>	$U_{11}$	$U_{22}$	$U_{33}$	$U_{12}$	$U_{13}$	$U_{23}$
Pb1	0.0149(6)	0.0141(6)	0.0078(6)	0.0000	0.0003(5)	0.0000
Pb2	0.0135(6)	0.0150(6)	0.0097(6)	0.0000	0.0004(5)	0.0000
Pb3	0.0146(6)	0.0142(6)	0.0084(6)	0.0000	0.0007(5)	0.0000
Pb4	0.0195(7)	0.0215(7)	0.0156(7)	0.0000	0.0004(5)	0.0000
Pt1	0.0204(7)	0.0145(6)	0.0098(7)	0.0000	-0.0011(5)	0.0000
Pt2	0.0196(7)	0.0153(6)	0.0110(7)	0.0000	-0.0016(5)	0.0000
Sr	0.0125(15)	0.0126(14)	0.0097(14)	0.0000	0.0028(12)	0.0000



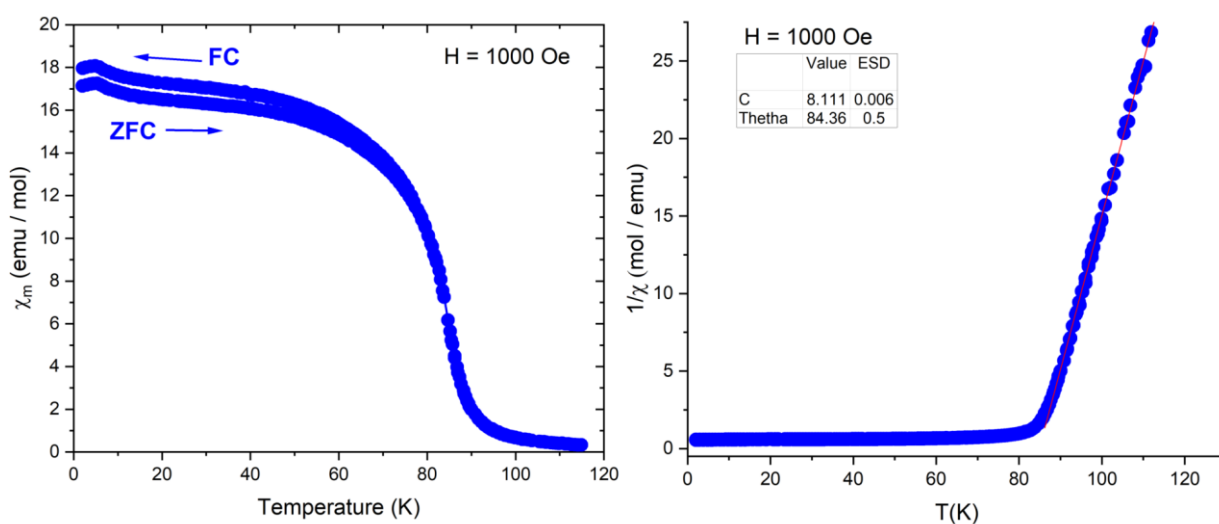
**Figure S2.** Observed, calculated, and difference Rietveld plots for the sample  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$ .

**Table S3.** Data collection and Rietveld analysis parameters, and final residuals for  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$

Compound	$\text{Eu}_2\text{Pt}_3\text{Pb}_5$
T, K	293
Space group	$Cmc2_1$
$a$ , Å	4.6430(9)
$b$ , Å	27.204(5)
$c$ , Å	7.562(2)
$V$ , Å <sup>3</sup>	962.1(3)
$Z$	4
Range in $2\theta$ , °	15-85
$\rho$ , g cm <sup>-3</sup>	13.291(5)
$R_{\text{Bragg}}$	0.022
$R_p$	0.046
$R_{\text{wp}}$	0.060
GoF	1.45

**Table S4.** Fractional atomic coordinates and isotropic displacement parameters for  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$  (polycrystalline sample).

Atom	Wickoff site	x/a	y/b	z/c	$B_{\text{eq}}$
Pt1	4a	0.5	0.5439(7)	1	0.75(1)
Pb1	4a	0	0.7139(6)	0	0.702(7)
Eu1	4a	0.5	0.6699(8)	1	0.805(5)
Eu2	4a	1	0.5195(6)	0	0.742(7)
Pb2	4a	1	0.7009(5)	0	0.955(8)
Pb3	4a	0.5	0.4524(7)	1	0.734(8)
Pt2	4a	0.5	0.7169(5)	0	0.734(8)
Pb4	4a	0.5	0.6174(4)	0	0.765(9)
Pt3	4a	1	0.6012(6)	0	0.773(8)
Pb5	4a	1	0.6001(6)	1	0.710(6)



**Figure S3.** The temperature dependences of molar magnetic susceptibility (left) of  $\text{Eu}_2\text{Pt}_3\text{Pb}_5$  and inverse magnetic susceptibility (right) in the magnetic field of 1000 Oe. The parameters of the Curie-Weiss fit are plotted in the inset table in the inverse magnetization plot.