

# Search Overview

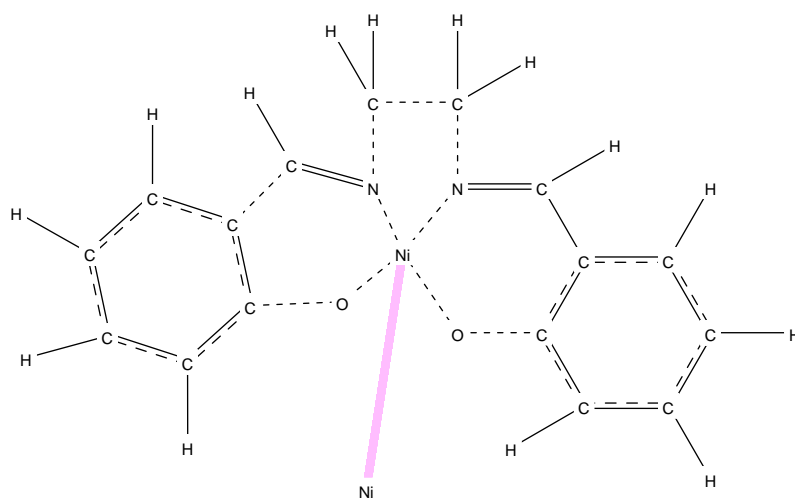
**Search:** search7  
**Date/Time done:** Wed Jan 30 14:32:42 2019  
**Database(s):** CSD version 5.40 (November 2018)  
**Restriction Info:** No refcode restrictions applied  
**Filters:** None  
**Percentage Completed:** 100%  
**Number of Hits:** 28

*Single query used. Search found structures that:*

match

**Query 1**

**Query 1**



# Search: search7 (Wed Jan 30 14:32:42 2019): Hits 1-4

## BEXKUK

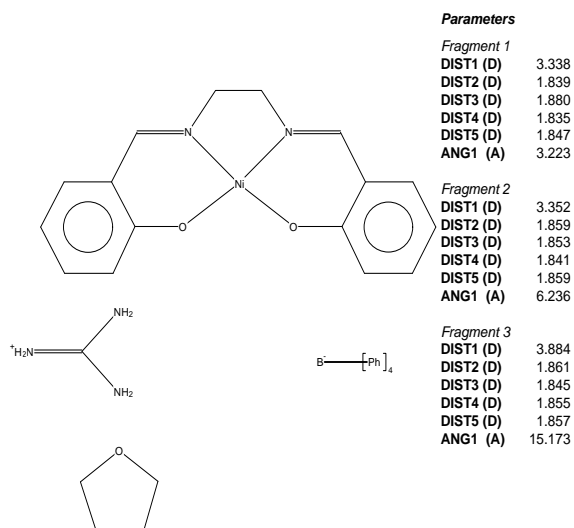
**Reference:** A.Giacomelli, C.Floriani, G.Perego (1982)  
*Chem.Commun.* ,650

**Formula:**  $3(C_{16}H_{14}N_2Ni_1O_2) \cdot C_1H_6N_3^{1+} \cdot C_{24}H_{20}B_1^{1-} \cdot 2(C_4H_8O_1)$

**Compound Name:** tris(N,N'-Ethylene-bis(salicylideneaminato))-nickel(ii) guanidinium tetraphenylborate tetrahydrofuran solvate

**Space Group:** P-1 **Cell:** a 16.496(8) b 19.228(8) c 15.226(4)  
**Space Group No.:** 2 **(Å,°)** α 94.88(3) β 102.83(3) γ 123.81(3)

**R-Factor (%):** 5.80 **Temperature(K):** 295 **Density(g/cm³):** 1.316



## COXNIN

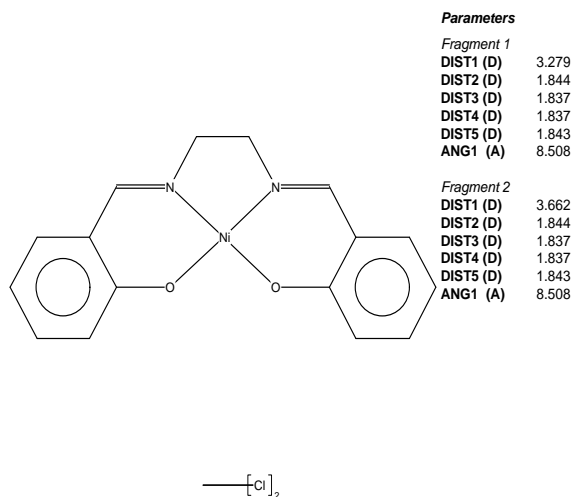
**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.* ,9,1194

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_1H_2Cl_2$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) dichloromethane solvate

**Space Group:** P21/n **Cell:** a 13.155(0) b 6.802(0) c 19.185(0)  
**Space Group No.:** 14 **(Å,°)** α 90.00 β 91.00(0) γ 90.00

**R-Factor (%):** 3.97 **Temperature(K):** 150 **Density(g/cm³):** 1.586



## COXNOT

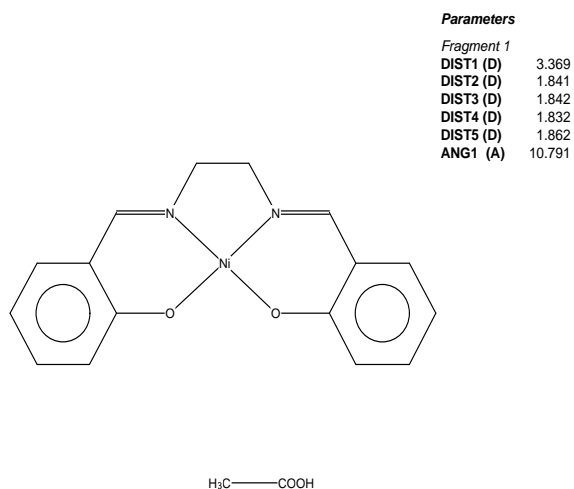
**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.* ,9,1194

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_2H_4O_2$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) acetic acid solvate

**Space Group:** P-1 **Cell:** a 11.494(0) b 7.362(0) c 10.015(0)  
**Space Group No.:** 2 **(Å,°)** α 93.81(0) β 102.23(0) γ 100.01(0)

**R-Factor (%):** 4.13 **Temperature(K):** 110 **Density(g/cm³):** 1.577



## COXNUZ

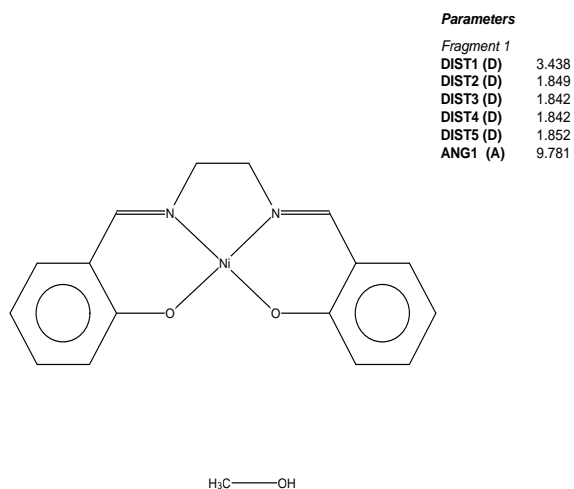
**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.* ,9,1194

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot 1.5(C_1H_4O_1)$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) methanol solvate

**Space Group:** Pbcn **Cell:** a 18.344(0) b 6.792(0) c 26.761(0)  
**Space Group No.:** 60 **(Å,°)** α 90.00 β 90.00 γ 90.00

**R-Factor (%):** 2.78 **Temperature(K):** 150 **Density(g/cm³):** 1.486



# Search: search7 (Wed Jan 30 14:32:42 2019): Hits 5-8

## FENRIA

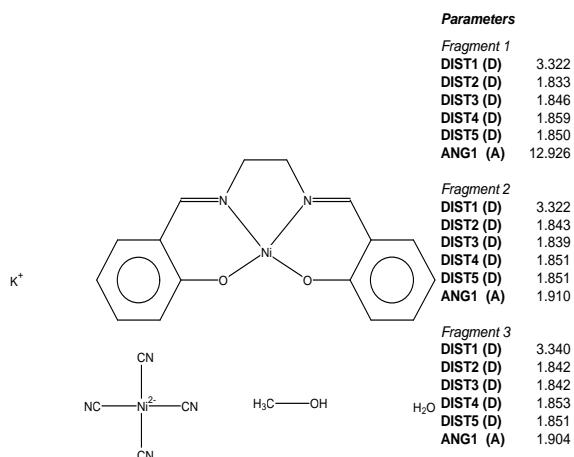
**Reference:** S.Skovgaard, A.D.Bond, C.J.McKenzie (2005) *Acta Crystallogr., Sect.E:Struct.Rep.Online* ,61,m135

**Formula:**  $2(K_1^{1+})_6(C_{16}H_{14}N_2Ni_1O_2)_4C_4N_4Ni_1^{2-} \cdot 6(C_1H_4O_1)_2(H_2O)_1$

**Compound Name:** Dipotassium hexakis((2,2'-(ethane-1,2-diylobis(nitrilomethylidyne)) diphenolato)-nickel(ii)) tetracyano-nickel(ii) methanol solvate dihydrate

**Space Group:** Pbc<sub>a</sub>    **Cell:**    *a* 14.321(1)    *b* 24.855(2)    *c* 29.709(3)  
**Space Group No.:** 61    (*Å*, °)     $\alpha$  90.00     $\beta$  90.00     $\gamma$  90.00

**R-Factor (%):** 5.47    **Temperature(K):** 180    **Density(g/cm<sup>3</sup>):** 1.520



## FEYWEM

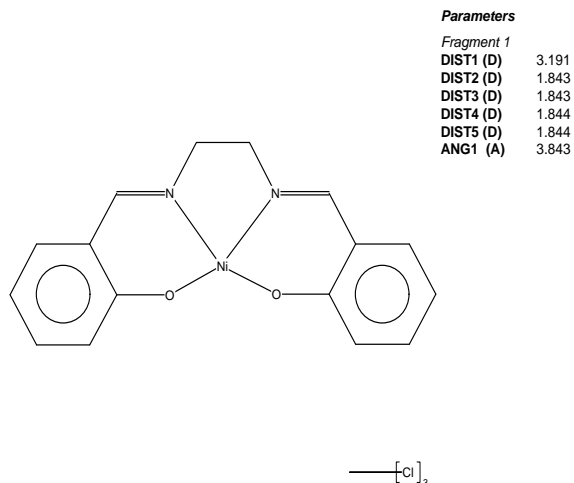
**Reference:** Xiao-Li Cheng, Shan Gao, Li-Hua Huo, S.W.Ng (2005) *Acta Crystallogr., Sect.E:Struct.Rep.Online* ,61,m385

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_1H_1Cl_3$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) chloroform solvate

**Space Group:** Pnmm    **Cell:**    *a* 6.997(1)    *b* 14.221(3)    *c* 18.355(4)  
**Space Group No.:** 58    (*Å*, °)     $\alpha$  90.00     $\beta$  90.00     $\gamma$  90.00

**R-Factor (%):** 4.60    **Temperature(K):** 295    **Density(g/cm<sup>3</sup>):** 1.616



## FEYWEM01

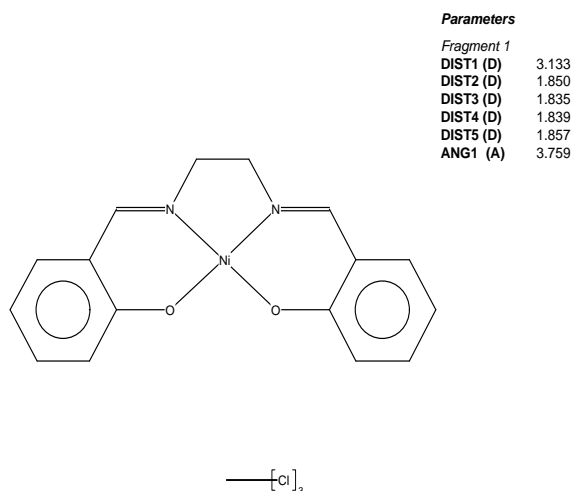
**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.* ,9,1194

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_1H_1Cl_3$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) chloroform solvate

**Space Group:** P21/n    **Cell:**    *a* 13.863(0)    *b* 6.949(0)    *c* 18.093(0)  
**Space Group No.:** 14    (*Å*, °)     $\alpha$  90.00     $\beta$  90.11(0)     $\gamma$  90.00

**R-Factor (%):** 3.70    **Temperature(K):** 110    **Density(g/cm<sup>3</sup>):** 1.694



## FEYWEM02

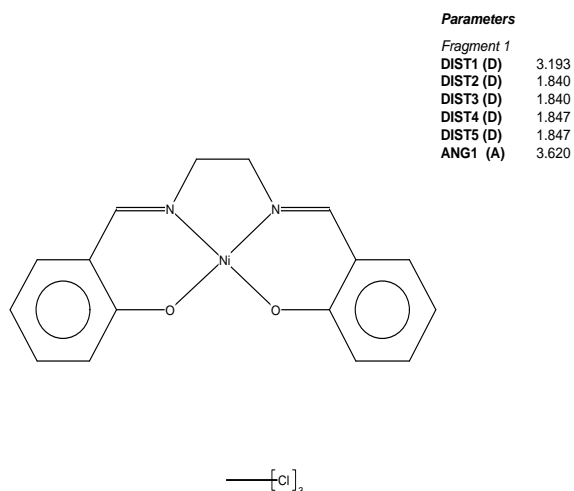
**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.* ,9,1194

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_1H_1Cl_3$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(ii) chloroform solvate

**Space Group:** Pnmm    **Cell:**    *a* 14.227(0)    *b* 6.995(0)    *c* 18.335(0)  
**Space Group No.:** 58    (*Å*, °)     $\alpha$  90.00     $\beta$  90.00     $\gamma$  90.00

**R-Factor (%):** 4.66    **Temperature(K):** 295    **Density(g/cm<sup>3</sup>):** 1.618



# Search: search7 (Wed Jan 30 14:32:42 2019): Hits 9-12

## JIMNOJ

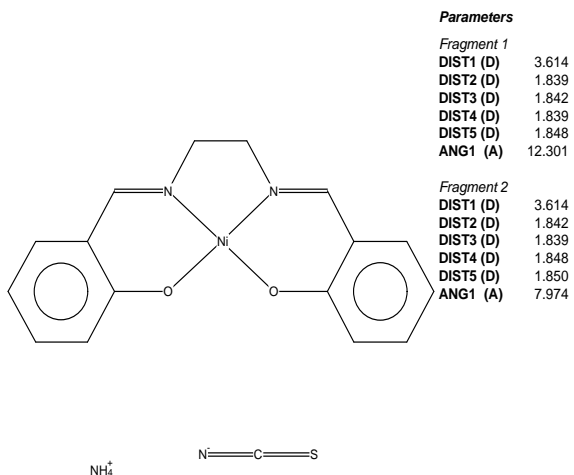
**Reference:** Xun Feng, Zhong-Xiang Du, Bao-Kian Ye, Fu-Na Cui (2007) *Jiegou Huaxue(Chin.)(Chin.J.Struct.Chem.)*, **26**,1033

**Formula:**  $2(C_{16}H_{14}N_2Ni_1O_2) \cdot H_4N_1^+ \cdot C_1N_1S_1^{1-}$

**Compound Name:** Ammonium bis(N,N'-ethylene-bis(salicylideneiminato)-nickel(ii) isothiocyanate

**Space Group:** Pbcu **Cell:** a 16.873(1) b 19.005(1) c 20.058(1)  
**Space Group No.:** 61 **(Å, °)** α 90.00 β 90.00 γ 90.00

**R-Factor (%):** 3.94 **Temperature(K):** 291 **Density(g/cm<sup>3</sup>):** 1.500



## NORFOR

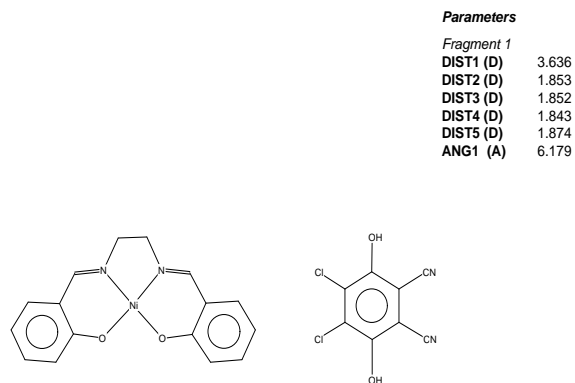
**Reference:** Ali Hossein Kianfar, M.Dostani, W.A.K.Mahmood (2015) *Polyhedron*, **85**,488

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2 \cdot C_8H_2Cl_2N_2O_2$

**Compound Name:** (2,2'-(ethane-1,2-diylbis((nitriolo)methylidene)diphenolato)-nickel(ii) 4,5-dichloro-3,6-dihydroxyphthalonitrile

**Space Group:** P-1 **Cell:** a 8.566(0) b 11.156(0) c 13.913(1)  
**Space Group No.:** 2 **(Å, °)** α 111.97(0) β 103.45(0) γ 92.79(0)

**R-Factor (%):** 4.21 **Temperature(K):** 100 **Density(g/cm<sup>3</sup>):** 1.552



## SAENNI

**Reference:** L.M.Shkol'nikova, E.M.Yumal', E.A.Shugam, V.A.Voblikova (1970) *Zh.Strukt.Khim.(Russ.)(J.Struct.Chem.)*, **11**,886

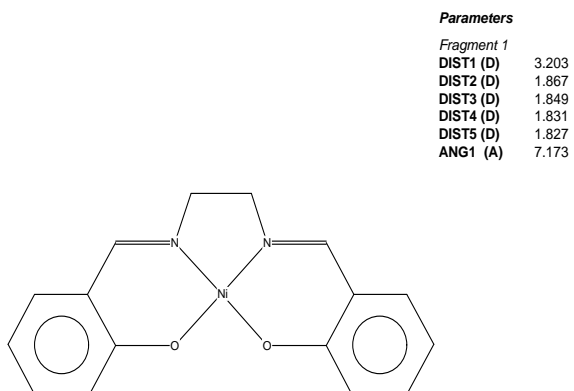
**Formula:**  $C_{16}H_{14}N_2Ni_1O_2$

**Compound Name:** (N,N'-bis(Salicylal)-ethylenedi-iminato)-nickel(ii)

**Synonym:** (2,2'-(ethane-1,2-diylbis((azanylylidene)methylidene)bis(phenolato))-nickel(ii)

**Space Group:** Pbcu **Cell:** a 13.818(8) b 26.132(9) c 7.483(3)  
**Space Group No.:** 61 **(Å, °)** α 90.00 β 90.00 γ 90.00

**R-Factor (%):** 12.70 **Temperature(K):** 295 **Density(g/cm<sup>3</sup>):** 1.597



## SAENNI01

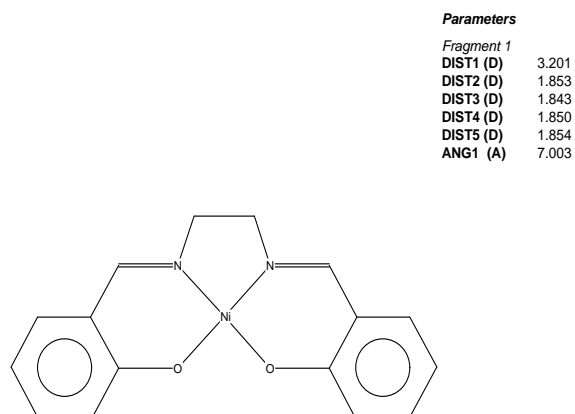
**Reference:** A.G.Manfredotti, C.Guastini (1983) *Acta Crystallogr.,Sect.C:Cryst.Struct Commun.*, **39**,863

**Formula:**  $C_{16}H_{14}N_2Ni_1O_2$

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato)-nickel(ii)

**Space Group:** Pbcu **Cell:** a 13.831(3) b 26.155(5) c 7.482(2)  
**Space Group No.:** 61 **(Å, °)** α 90.00 β 90.00 γ 90.00

**R-Factor (%):** 4.20 **Temperature(K):** 295 **Density(g/cm<sup>3</sup>):** 1.595



# Search: search7 (Wed Jan 30 14:32:42 2019): Hits 13-16

## SAENNI02

**Reference:** E.F.DiMauro, M.C.Kozlowski (2002) *Organometallics*, **21**, 1454

**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>

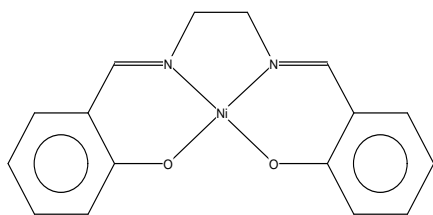
**Compound Name:** (N,N'-Ethylene-bis(salicylideneaminato-N,O))-nickel(II)

**Space Group:** Pbc<sub>a</sub>    **Cell:**    **a** 7.429(0)    **b** 13.740(0)    **c** 26.069(0)  
**Space Group No.:** 61    **(Å, °)**    α 90.00    β 90.00    γ 90.00

**R-Factor (%):** 6.29    **Temperature(K):** 200    **Density(g/cm<sup>3</sup>):** 1.623

### Parameters

Fragment 1  
**DIST1 (D)** 3.180  
**DIST2 (D)** 1.850  
**DIST3 (D)** 1.850  
**DIST4 (D)** 1.850  
**DIST5 (D)** 1.851  
**ANG1 (A)** 6.946



## SAENNI03

**Reference:** M.Kondo, K.Nabari, T.Horiba, Y.Irie, Md K.Kabir, R.P.Sarker, E.Shimizu, Y.Shimizu, Y.Fuwa (2003) *Inorg.Chem.Commun.*, **6**,154

**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>

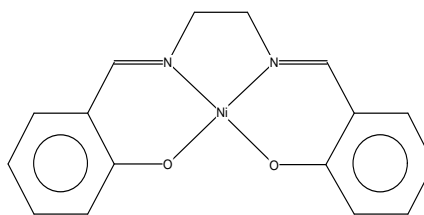
**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(II)

**Space Group:** Pbc<sub>a</sub>    **Cell:**    **a** 7.484(3)    **b** 13.822(6)    **c** 26.150(10)  
**Space Group No.:** 61    **(Å, °)**    α 90.00    β 90.00    γ 90.00

**R-Factor (%):** 4.76    **Temperature(K):** 293    **Density(g/cm<sup>3</sup>):** 1.596

### Parameters

Fragment 1  
**DIST1 (D)** 3.201  
**DIST2 (D)** 1.846  
**DIST3 (D)** 1.853  
**DIST4 (D)** 1.852  
**DIST5 (D)** 1.851  
**ANG1 (A)** 6.759



## SAENNI05

**Reference:** M.A.Siegler, M.Lutz (2009) *Cryst.Growth Des.*, **9**,1194

**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>

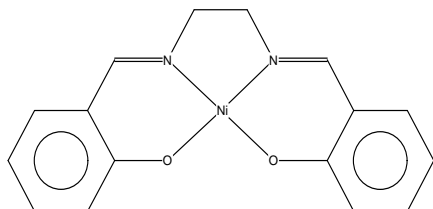
**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato))-nickel(II)

**Space Group:** Pcab    **Cell:**    **a** 13.684(0)    **b** 7.377(0)    **c** 25.968(0)  
**Space Group No.:** 61    **(Å, °)**    α 90.00    β 90.00    γ 90.00

**R-Factor (%):** 1.86    **Temperature(K):** 100    **Density(g/cm<sup>3</sup>):** 1.647

### Parameters

Fragment 1  
**DIST1 (D)** 3.160  
**DIST2 (D)** 1.852  
**DIST3 (D)** 1.849  
**DIST4 (D)** 1.849  
**DIST5 (D)** 1.854  
**ANG1 (A)** 6.913



## SAENNI06

**Reference:** Y.Ding, F.Wang, Z.J.Ku, L.S.Wang, Q.R.Wang (2009) *Koord.Khim.(Russ.)(Coord.Chem.)*, **35**,366

**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>

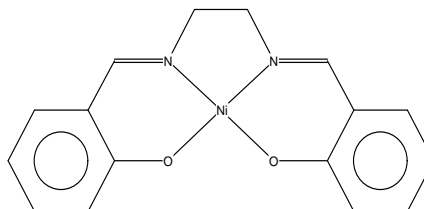
**Compound Name:** (Ethylene-1,2-bis(salicylideneaminato))-nickel(II)

**Space Group:** Pbc<sub>a</sub>    **Cell:**    **a** 7.472(0)    **b** 13.805(0)    **c** 26.098(1)  
**Space Group No.:** 61    **(Å, °)**    α 90.00    β 90.00    γ 90.00

**R-Factor (%):** 2.38    **Temperature(K):** 295    **Density(g/cm<sup>3</sup>):** 1.604

### Parameters

Fragment 1  
**DIST1 (D)** 3.196  
**DIST2 (D)** 1.845  
**DIST3 (D)** 1.850  
**DIST4 (D)** 1.852  
**DIST5 (D)** 1.845  
**ANG1 (A)** 6.699



# Search: search7 (Wed Jan 30 14:32:42 2019): Hits 17-20

## SAENNI07

**Reference:** J.G.Malecki (2012)  
CSD Communication(Private Communication) ,

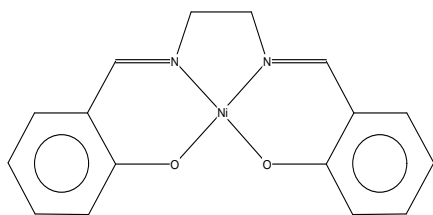
**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>

**Compound Name:** (N,N'-Ethylenebis(salicylideneiminato))-nickel(ii)

**Space Group:** Pbc<sub>a</sub> **Cell:** a 7.474(1) b 13.808(3) c 26.143(5)  
**Space Group No.:** 61 **(Å, °)** α 90.00 β 90.00 γ 90.00  
**R-Factor (%):** 3.95 **Temperature(K):** 295 **Density(g/cm<sup>3</sup>):** 1.600

### Parameters

**Fragment 1**  
**DIST1 (D)** 3.198  
**DIST2 (D)** 1.852  
**DIST3 (D)** 1.845  
**DIST4 (D)** 1.852  
**DIST5 (D)** 1.850  
**ANG1 (A)** 6.596



## TUNREB

**Reference:** R.L.De, M.Mandal, L.Roy, J.Mukherjee, R.K.De (2014)  
J.Indian Chem.Soc. ,91,2165

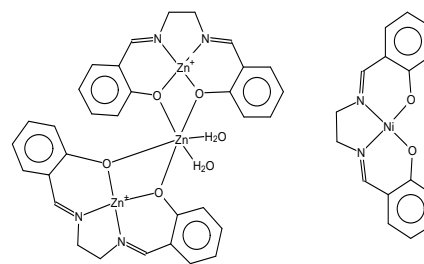
**Formula:** C<sub>32</sub> H<sub>32</sub> N<sub>4</sub> O<sub>6</sub> Zn<sub>3</sub> 2+(C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>).2(Cl<sub>1</sub> O<sub>4</sub> 1-).0.5(H<sub>2</sub> O<sub>1</sub>)

**Compound Name:** bis(μ<sub>2</sub>-2,2'-(Ethane-1,2-diybis(nitriolomethylidene)diphenolato)-diaqua-tri-zinc (2,2'-(ethane-1,2-diybis(nitriolomethylidene)diphenolato)-nickel diperchlorate hemihydrate

**Space Group:** C2/c **Cell:** a 28.527(2) b 10.175(0) c 24.161(2)  
**Space Group No.:** 15 **(Å, °)** α 90.00 β 111.89(0) γ 90.00  
**R-Factor (%):** 6.48 **Temperature(K):** 293 **Density(g/cm<sup>3</sup>):** 1.656

### Parameters

**Fragment 1**  
**DIST1 (D)** 3.281  
**DIST2 (D)** 1.837  
**DIST3 (D)** 1.836  
**DIST4 (D)** 1.853  
**DIST5 (D)** 1.858  
**ANG1 (A)** 1.744



H<sub>2</sub>O

ClO<sub>4</sub><sup>-</sup>

## UMIBAT

**Reference:** M.Lutz (2003) Acta Crystallogr.,Sect.E:Struct.Rep.Online ,  
59,m950

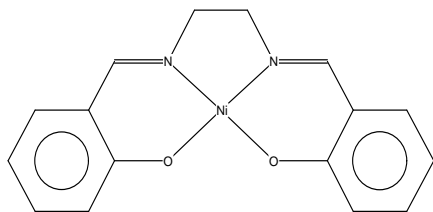
**Formula:** C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>.C<sub>3</sub> H<sub>7</sub> N<sub>1</sub> O<sub>1</sub>

**Compound Name:** (N,N'-Ethylene-bis(salicylideneiminato-N,O))-nickel(ii) dimethylformamide solvate

**Space Group:** P2<sub>1</sub>/c **Cell:** a 13.387(0) b 6.669(0) c 22.733(0)  
**Space Group No.:** 14 **(Å, °)** α 90.00 β 118.04(0) γ 90.00  
**R-Factor (%):** 2.70 **Temperature(K):** 150 **Density(g/cm<sup>3</sup>):** 1.476

### Parameters

**Fragment 1**  
**DIST1 (D)** 3.390  
**DIST2 (D)** 1.848  
**DIST3 (D)** 1.845  
**DIST4 (D)** 1.845  
**DIST5 (D)** 1.850  
**ANG1 (A)** 7.132



OHC—NMe<sub>2</sub>

## VUYJAB

**Reference:** G.Assey, Y.Gultneh, R.J.Butcher (2010)  
Acta Crystallogr.,Sect.E:Struct.Rep.Online ,66,m654

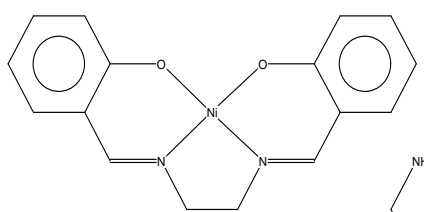
**Formula:** C<sub>2</sub> H<sub>10</sub> N<sub>2</sub> 2+.4(C<sub>16</sub> H<sub>14</sub> N<sub>2</sub> Ni<sub>1</sub> O<sub>2</sub>).2(Cl<sub>1</sub> O<sub>4</sub> 1-).2(C<sub>3</sub> H<sub>7</sub> N<sub>1</sub> O<sub>1</sub>)

**Compound Name:** Ethylenediammonium tetrakis((2,2'-(ethane-1,2-diybis(nitriolomethylidene)diphenolato)-nickel(ii)) diperchlorate N,N-dimethylformamide solvate

**Space Group:** P-1 **Cell:** a 15.021(1) b 15.049(1) c 18.271(0)  
**Space Group No.:** 2 **(Å, °)** α 85.99(0) β 86.51(0) γ 62.96(0)  
**R-Factor (%):** 3.89 **Temperature(K):** 200 **Density(g/cm<sup>3</sup>):** 1.546

### Parameters

**Fragment 1**  
**DIST1 (D)** 3.454  
**DIST2 (D)** 1.844  
**DIST3 (D)** 1.850  
**DIST4 (D)** 1.838  
**DIST5 (D)** 1.852  
**ANG1 (A)** 11.970



Me<sub>2</sub>N—CHO

NH<sub>3</sub><sup>+</sup>  
ClO<sub>4</sub><sup>-</sup>  
H<sub>3</sub>N<sup>+</sup>

**Fragment 2**  
**DIST1 (D)** 3.454  
**DIST2 (D)** 1.848  
**DIST3 (D)** 1.837  
**DIST4 (D)** 1.849  
**DIST5 (D)** 1.837  
**ANG1 (A)** 2.806

**Fragment 3**  
**DIST1 (D)** 3.644  
**DIST2 (D)** 1.846  
**DIST3 (D)** 1.845  
**DIST4 (D)** 1.841  
**DIST5 (D)** 1.856  
**ANG1 (A)** 10.723

**Fragment 4**  
**DIST1 (D)** 3.644  
**DIST2 (D)** 1.850  
**DIST3 (D)** 1.842  
**DIST4 (D)** 1.856  
**DIST5 (D)** 1.845  
**ANG1 (A)** 13.120