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Supporting Information

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1. General and Materials

General: All reactions were carried out under an atmosphere of nitrogen using the standard Schlenk techniques, unless otherwise noted. ¹H NMR and ¹³C NMR spectra were recorded at room temperature in CDCl₃, CD₃OD, DMSO-d₆ on 400 MHz instrument with tetramethylsilane (TMS) as internal standard. Enantiomeric excess was determined by HPLC analysis, using chiral column described below in detail. Optical rotations were measured by polarimeter. Flash column chromatography was performed on silica gel (200-300 mesh). All reactions were monitored by TLC analysis.

Materials: Commercially available reagents were used throughout without further purification. The anhydrous solvents for asymmetric hydrogenation were also purchased without the further purification.

2. Synthesis of Quinazolinone Derivatives

Quinazolinone derivatives **1** can be conveniently synthesized according to the known literature procedure.¹ Among them, the quinazolinones **1a**,¹ **1d**,² **1g**,³ **1i**,⁴ **1j**,⁵ **1k**,⁶ **1l**,⁷ **1m**,⁸ **1n**⁹ and **1o**¹⁰ are the known compounds.



General procedure: the Grignand reagent was prepared by reaction of magnesium (288 mg, 12 mmol) with the corresponding aryl bromide (15.6 mmol) in dry tetrahydrofuran under reflux. Then, the 2-aminobenzonitrile (708 mg, 6.0 mmol) in dry tetrahydrofuran (8 mL) was added dropwise under reflux. After a refluxed period (2 h), the mixture was cooled to 0 $^{\circ}$ C, methyl chloroformate (977 mg, 9.0 mmol) was added dropwise, and the solution was refluxed for 14 h. The mixture was cooled to room temperature and poured into the hydrochloric acid solution (2 M), then neutralized with 10% sodium bicarbonate solution and extracted with dichloromethane. The combined organic layer was dried over anhydrous sodium sulfate, concentrated in *vacuo*. The residue was further purified by flash column chromatography using dichloromethane/methanol as eluent to afford the desired quinazolinones **1**.

4-*o***-Tolylquinazolin-2(1***H***)-one (1b): 1.066 g, 75% yield, white solid, mp: 262-263 °C, new compound, R_f = 0.30 (dichloromethane/methanol = 15/1); ¹H NMR (400 MHz, CDCl₃) \delta 13.05 (s,**



1H), 7.73–7.68 (m, 1H), 7.59 (d, J = 8.2 Hz, 1H), 7.45–7.41 (m, 2H), 7.36–7.33 (m, 3H), 7.20–7.16 (m, 1H), 2.26 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 179.4, 159.3, 143.6, 137.0, 136.5, 136.4, 131.4, 130.4, 129.5, 129.4, 126.4, 124.1, 117.5, 117.2, 20.6; HRMS (ESI) *m*/*z* Calculated for C₁₅H₁₃N₂O [M+H]⁺ 237.1022,

4-*m***-Tolylquinazolin-2(1***H***)-one (1c): 1.134 g, 80% yield, white solid, mp: 232-233 °C, new compound, R_f = 0.30 (dichloromethane/methanol = 15/1); ¹H NMR (400 MHz, CDCl₃) \delta 7.90 (d,**



J = 8.2 Hz, 1H), 7.73–7.58 (m, 4H), 7.45–7.39 (m, 2H), 7.30–7.16 (m, 1H),, 2.47 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 176.6, 158.2, 143.2, 138.2, 136.3, 135.1, 131.4, 130.3, 128.7, 128.0, 126.8, 122.9, 116.6, 115.3, 21.3; HRMS (ESI) *m*/*z* Calculated for C₁₅H₁₃N₂O [M+H]⁺ 237.1022, found 237.1024. **4-(3,5-Dimethylphenyl)quinazolin-2(1***H***)-one (1e):** 0.945 g (4.0 mmol scale), 95% yield, white solid, mp: 290-291 °C, new compound, $R_f = 0.30$ (dichloromethane/methanol = 15/1); ¹H



NMR (400 MHz, CDCl₃) δ 7.91 (d, J = 8.2 Hz, 1H), 7.73–7.62 (m, 2H), 7.43 (s, 2H), 7.27–7.21 (m, 2H), 2.43 (s, 6H); ¹³C NMR (100 MHz, CDCl₃) δ 177.8, 159.2, 144.2, 138.9, 137.3, 136.0, 133.3, 129.8, 128.5, 123.8, 117.6, 116.3, 22.1; HRMS (ESI) m/z Calculated for C₁₆H₁₅N₂O [M+H]⁺ 251.1179, found 251.1181.

4-(3-Methoxyphenyl)quinazolin-2(1*H***)-one (1f):** 0.920 g (4.0 mmol scale), 92% yield, white solid, mp: 254-255 °C, new compound, $R_f = 0.32$ (dichloromethane/methanol = 15/1); ¹H NMR



(400 MHz, DMSO-d₆) δ 7.73 (t, J = 7.8 Hz, 1H), 7.65 (d, J = 8.2 Hz, 1H), 7.48 (t, J = 7.8 Hz, 1H), 7.37 (d, J = 8.4 Hz, 1H), 7.22–7.15 (m, 4H), 3.82 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 176.1, 160.4, 144.7, 139.1, 136.4, 130.9, 129.6, 123.6, 122.6, 117.4, 116.8, 115.5, 115.5, 110.8, 56.6;

HRMS (ESI) m/z Calculated for C₁₅H₁₃N₂O₂ [M+H]⁺ 253.0972, found 253.0974.

4-(3,5-Dimethoxyphenyl)quinazolin-2(1*H***)-one (1h):** 0.857 g (4.0 mmol scale), 76% yield, white solid, mp: 255-256 °C, new compound, $R_f = 0.30$ (dichloromethane/methanol = 15/1); ¹H



NMR (400 MHz, CDCl₃) δ 7.94 (d, J = 8.2 Hz, 1H), 7.72 (t, J = 7.6 Hz, 1H), 7.61 (d, J = 8.2 Hz, 1H), 7.26 (d, J = 6.3 Hz, 2H), 6.92 (d, J = 1.4 Hz, 2H), 6.67 (s, 1H), 3.87 (s, 6H); ¹³C NMR (100 MHz, CDCl₃) δ 177.4, 161.5, 159.0, 144.1, 139.1, 136.2, 129.7, 123.9, 117.5, 116.2, 108.6, 103.9, 56.5; HRMS (ESI) m/z Calculated for C₁₆H₁₅N₂O₃ [M+H]⁺ 283.1077,

found 283.1079.

3. General Procedure for Asymmetric Hydrogenation of Quinazolinones



A mixture of $[Ir(cod)Cl]_2$ (1.3 mg, 0.002 mmol) and (*R*)-SegPhos (2.8 mg, 0.0044 mmol) in tetrahydrofuran (1.0 mL) was stirred at room temperature for 10 min in a glovebox, then BCDMH (4.8 mg, 0.02 mmol) and substrates **1** (0.2 mmol) together with tetrahydrofuran (2.0 mL) were added and the mixture was stirred for a further 10 min. The hydrogenation was performed at 25 °C under hydrogen gas (600 psi) in a stainless steel autoclave for 24 h. After carefully releasing the hydrogen gas, saturated aqueous sodium bicarbonate (3.0 mL) was added into the mixture and stirred for 10-15 min. The mixture was extracted with dichloromethane three times and the combined organic extract was dried over anhydrous sodium sulfate. After filtration, the filtrate was concentrated in *vacuo* and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate (or dichloromethane/methanol) to give the desired products **2**.

(S)-(-)-4-Phenyl-3,4-dihydroquinazolin-2(1*H*)-one (2a): 41 mg, 91% yield, white solid, the known compound,¹⁰ 98% ee, $[\alpha]_{D}^{20} = -110.4$ (*c* 0.92, MeOH), $R_f = 0.31$ (dichloromethane/

HN NH (m HN 144 En

methanol = 15/1); ¹H NMR (400 MHz, CDCl₃) δ 7.35–7.28 (m, 5H), 7.14–7.10 (m, 1H), 6.88–6.79 (m, 3H), 5.64 (s, 1H); ¹³C NMR (100 MHz, CD₃OD) δ 155.4, 144.2, 136.0, 128.4, 127.9, 127.5, 126.7, 126.5, 122.0, 121.5, 114.0, 57.8; Enantiomeric excess was determined by HPLC (OD-H column, *n*-Hexane/

i-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 16.9$ min (major), $t_2 = 20.7$ min.

(-)-4-o-Tolyl-3,4-dihydroquinazolin-2(1H)-one (2b): 46 mg, 97% yield, white solid, mp: 117-118 °C, new compound, 92% ee, $[\alpha]^{20}_{D} = -74.2$ (c 0.26, MeOH), $R_f = 0.35$ (dichloromethane



HN

/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.24 (s, 1H), 7.21–7.09 (m, 6H), 6.85–6.82 (m, 1H), 6.79–6.75 (m, 1H), 6.70 (d, *J* = 7.2 Hz, 1H), 5.81 (d, *J* = 1.8 Hz, 1H), 2.39 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 153.7, 143.0, 137.8, 135.5, 131.2, 128.6, 128.3, 127.9, 126.9, 126.8, 121.7, 121.6, 114.3, 54.7, 19.5;

Enantiomeric excess was determined by HPLC (OD-H column, n-Hexane/i-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 12.6$ min (major), $t_2 = 14.7$ min; HRMS (ESI) m/z Calculated for C₁₅H₁₅N₂O [M+H]⁺ 239.1179, found 239.1177.

(-)-4-m-Tolyl-3,4-dihydroquinazolin-2(1H)-one (2c): 42 mg, 90% yield, white solid, mp: 210-211 °C, new compound, 96% ee, $[\alpha]_{D}^{20} = -78.3$ (*c* 0.66, MeOH), $R_f = 0.35$ (dichloromethane

/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.21 (s, 1H), 7.38 (s, 1H), 7.20 (t, J = 7.5 Hz, 1H), 7.10–7.00 (m, 5H), 6.80 (t, J = 6.8 Hz, 2H), 5.47 (s, NH 1H), 2.26 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 154.2, 145.5, 138.1, 137.4, 128.9, 128.4, 128.3, 127.2, 123.9, 122.1, 121.5, 114.3, 57.2, 21.6; Enantiomeric excess was determined by HPLC (OD-H column, n-Hexane/*i*-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 16.5 \text{ min (major)}, t_2 = 19.7 \text{ min; HRMS (ESI)}$

m/z Calculated for C₁₅H₁₅N₂O [M+H]⁺ 239.1179, found 239.1177.

(-)-4-p-Tolyl-3,4-dihydroquinazolin-2(1H)-one (2d): 45 mg, 97% yield, white solid, known compound,² 95% ee, $[\alpha]_{D}^{20}$ = -135.8 (c 0.33, MeOH), R_f = 0.30 (dichloromethane/ methanol =



15/1); ¹H NMR (400 MHz, CDCl₃) δ 8.96 (s, 1H), 7.26–7.22 (m, 2H), 7.15– 7.07 (m, 3H), 6.86–6.78 (m, 3H), 5.93 (s, 1H), 5.60 (s, 1H), 2.31 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 155.2, 140.1, 138.0, 136.0, 129.6, 128.3, 127.1, 126.9, 122.3, 121.5, 114.6, 58.3, 21.1; Enantiomeric excess was determined by

chiral HPLC (OD-H column, *n*-Hexane/*i*-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30° C), t₁ = 18.5 min (major), t₂ = 22.6 min.

(-)-4-(3,5-Dimethylphenyl)-3,4-dihydroquinazolin-2(1H)-one (2e): 46 mg, 91% yield, white solid, mp: 200-201 °C, new compound, 96% ee, $[\alpha]^{20}_{D}$ = -140.8 (c 0.24, MeOH), R_f = 0.35



(dichloromethane/methanol = 15:1); ¹H NMR (400 MHz, CDCl₃) δ 7.81 (s, 1H), 7.17–7.13 (m, 1H), 6.96 (s, 3H), 6.89–6.82 (m, 2H), 6.76 (d, J = 7.8 Hz, 1H), 5.58 (s, 1H), 5.27 (s, 1H), 2.30 (s, 6H); ¹³C NMR (100 MHz, CDCl₃) δ 154.3, 142.7, 138.7, 135.7, 130.0, 128.4, 127.1, 125.0, 122.4, 121.4, 114.3, 58.8, 21.3; Enantiomeric excess was determined by HPLC (OD-H column,

n-Hexane /i-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), t₁ = 13.1 min (major), $t_2 = 16.8 \text{ min}; \text{HRMS (ESI)} m/z \text{ Calculated for } C_{16}H_{17}N_2O[M+H]^+ 253.1335, \text{ found } 253.1337.$

(-)-4-(3-Methoxyphenyl)-3,4-dihydroquinazolin-2(1H)-one (2f): 45 mg, 88% yield, white solid, mp: 175-176 °C, new compound, 95% ee, $[\alpha]^{20}_{D} = -70.4$ (c 0.82, MeOH), $R_f = 0.40$ (neat



ethyl acetate); ¹H NMR (400 MHz, DMSO-d₆) δ 9.26 (s, 1H), 7.44 (s, 1H), 7.24 (t, J = 7.8 Hz, 1H), 7.13–7.07 (m, 2H), 6.88–6.81 (m, 5H), 5.50 (s, 1H), 3.72 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 159.8, 154.2, 147.0, 137.4, 130.2, 128.3, 127.2, 122.0, 121.6, 118.8, 114.4, 112.8, 112.7, 57.0,

55.5; Enantiomeric excess was determined by HPLC (IC column, n-Hexane/i-PrOH = 80/20,

detector: 254 nm, flow rate: 0.80 mL/min, 30 °C), $t_1 = 15.7$ min, $t_2 = 25.8$ min (major); HRMS (ESI) m/z Calculated for C₁₅H₁₅N₂O₂ [M+H]⁺ 255.1128, found 255.1126.

(-)-4-(4-Methoxyphenyl)-3,4-dihydroquinazolin-2(1H)-one (2g): 50 mg, 98% yield, white solid, mp: 236-237 °C, new compound, 91% ee, $[\alpha]_{D}^{20}$ = -85.6 (c 0.55, MeOH), $R_f = 0.40$

(dichloromethane/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.24 NH (s, 1H), 7.38 (s, 1H), 7.21 (d, J = 8.6 Hz, 2H), 7.11 (t, J = 7.6 Hz, 1H), 7.00

> (d, J = 7.6 Hz, 1H), 6.89 (d, J = 8.6 Hz, 2H), 6.82 (t, J = 7.4 Hz, 2H), 5.48 OMe (d, J = 1.8 Hz, 1H), 3.72 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 159.0,

154.2, 137.6, 137.4, 128.2, 128.0, 127.2, 122.5, 121.5, 114.3, 56.6, 55.6; Enantiomeric excess was determined by HPLC (OD-3 column, n-Hexane/i-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 24.0$ min, $t_2 = 25.7$ min (major); HRMS (ESI) m/z Calculated for $C_{15}H_{15}N_2O_2$ [M+H]⁺ 255.1128, found 255.1133.

(-)-4-(3,5-Dimethoxyphenyl)-3,4-dihydroquinazolin-2(1H)-one (2h): 52 mg, 91% yield, white solid, mp: 115-116 °C, new compound, 95% ee, $[\alpha]^{20}_{D}$ = -85.5 (*c* 0.60, MeOH), R_f = 0.30



(dichloromethane/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.19 (s, 1H), 7.36 (s, 1H), 7.10-7.05 (m, 2H), 6.82-6.76 (m, 2H), 6.42 (d, J = 1.8 Hz, 2H), 6.36 (s, 1H), 5.40 (d, J = 1.8 Hz, 1H), 3.67 (s, 6H); ¹³C NMR (100 MHz, DMSO-d₆) δ 160.5, 153.7, 147.2, 136.9, 127.8, 126.7, 121.3, 121.0, 113.8, 104.5, 98.4, 56.6, 55.1; Enantiomeric excess was

determined by HPLC (IA column, n-Hexane/i-PrOH = 75/25, detector: 254 nm, flow rate: 0.90 mL/min, 30 °C), $t_1 = 12.0$ min (major), $t_2 = 21.0$ min; HRMS (ESI) m/z Calculated for $C_{16}H_{17}N_2O_3$ [M+H]⁺ 285.1234, found 285.1236.

(-)-4-(4-Chlorophenyl)-3,4-dihydroquinazolin-2(1H)-one (2i): 50 mg, 97% yield, white solid, the known compound,¹¹ 98% ee, $[\alpha]_{D}^{20} = -168.9$ (c 0.54, MeOH), $R_f = 0.42$ (dichloromethane/



methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.27 (s, 1H), 7.47 (s, 1H), 7.40 (d, J = 8.6 Hz, 2H), 7.31 (d, J = 8.4 Hz, 2H), 7.12 (t, J = 7.6 Hz, 1H), 7.05 (d, J = 7.4 Hz, 1H), 6.83 (t, J = 7.6 Hz, 2H), 5.56 (s, 1H); ¹³C NMR (100 MHz, DMSO-d₆) δ 154.1, 144.4, 137.4, 132.4, 129.0, 128.6, 128.5, 127.2, 121.7, 121.6, 114.5, 56.4; Enantiomeric excess was determined by HPLC (OD-3 column,

n-Hexane/*i*-PrOH = 90/10, detector: 254 nm, flow rate: 0.90 mL/min, 30 °C), $t_1 = 19.3$ min, $t_2 = 19.3$ 20.1 min (major).

(-)-4-(4-Fluorophenyl)-3,4-dihydroquinazolin-2(1H)-one (2j): 46 mg, 95% yield, white solid, the known compound [CAS: 1781596-98-2], 97% ee, $[\alpha]_{D}^{20} = -88.3$ (c 0.30, MeOH), $R_{f} = 0.38$



(dichloromethane/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.32 (s, 1H), 7.50 (s, 1H), 7.36–7.32 (m, 2H), 7.19–7.10 (m, 3H), 7.04 (d, J = 7.2 Hz, 1H), 6.86–6.82 (m, 2H), 5.58 (d, J = 2.4 Hz, 1H); ¹³C NMR (100 MHz, DMSO-d₆) δ 161.9 (d, J = 243.2 Hz), 154.2, 141.7, 141.6, 137.4, 128.8 (d, J =

8.3 Hz), 128.4, 127.2, 121.9, 121.7, 115.7 (d, J = 21.4 Hz), 114.5, 56.4; ¹⁹F NMR (376 MHz, DMSO-d₆) δ -115.36; Enantiomeric excess was determined by HPLC (OD-3 column, *n*-Hexane /i-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 15.4$ min, $t_2 = 16.6$ min (major).

(-)-6-Methyl-4-phenyl-3,4-dihydroquinazolin-2(1H)-one (2k): 45 mg, 95% yield, white solid, the known compound, ¹² 97% ee, $[\alpha]^{20}_{D} = -31.1$ (c 0.46, MeOH), $R_f = 0.25$ (dichloromethane/ methanol = 30/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.18 (s, 1H), 7.39–7.22 (m, 6H), 6.92 (d, J =



8.0 Hz, 1H), 6.85 (s, 1H), 6.73 (d, J = 8.0 Hz, 1H), 5.47 (s, 1H), 2.14 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 154.3, 145.6, 135.0, 130.3, 129.0, 128.8, 127.8, 127.5, 126.7, 121.9, 114.3, 57.3, 20.8; Enantiomeric excess was determined by HPLC (OD-H column, *n*-Hexane/*i*-PrOH = 95/05, detector: 254 nm, flow rate: 0.70 mL/ min, 30 °C), t₁ = 44.7 min (major), t₂ = 51.8 min.

(-)-6-Chloro-4-phenyl-3,4-dihydroquinazolin-2(1*H*)-one (2l): 50 mg, 97% yield, white solid, the known compound,¹⁰ 96% ee, $[\alpha]^{20}_{D}$ = -12.2 (*c* 0.49, MeOH), R_f = 0.23 (dichloromethane/



methanol = 25/1); ¹H NMR (400 MHz, DMSO-d₆) δ 9.40 (s, 1H), 7.54 (s, 1H), 7.37–7.25 (m, 5H), 7.18–7.04 (m, 2H), 6.83 (d, *J* = 8.4 Hz, 1H), 5.55 (d, *J* = 2.4 Hz, 1H); ¹³C NMR (100 MHz, DMSO-d₆) δ 153.9, 145.0, 136.5, 129.2, 128.2, 128.0, 126.8, 126.6, 125.0, 124.1, 116.0, 56.6; Enantiomeric excess was determined by HPLC (OD-H column, *n*-Hexane/*i*-PrOH = 90/10, detector: 254 nm,

flow rate: 0.70 mL/min, 30 °C), $t_1 = 21.4 \text{ min (major)}, t_2 = 23.9 \text{ min.}$

(-)-6,7-Dimethoxy-4-phenyl-3,4-dihydroquinazolin-2(1*H*)-one (2m): 48 mg, 85% yield, white solid, mp: 110- 111 °C, new compound, 93% ee, $[\alpha]_{D}^{20} = -53.5$ (*c* 0.54, MeOH), $R_f = 0.35$



(dichloromethane/methanol = 25/1); ¹H NMR (400 MHz, DMSO-d₆) δ 8.97 (s, 1H), 7.34–7.28 (m, 5H), 7.25–7.21 (m, 1H), 6.69 (s, 1H), 6.48 (s, 1H), 5.42 (d, *J* = 2.6 Hz, 1H), 3.68 (d, *J* = 4.0 Hz, 3H), 3.60 (s, 3H); ¹³C NMR (100 MHz, DMSO-d₆) δ 154.2, 149.2, 145.7, 143.9, 131.2, 129.0, 127.6, 126.6, 113.0, 111.6, 99.3, 57.0, 56.6, 55.9; Enantiomeric excess was deter-

mined by HPLC (IC column, *n*-Hexane/*i*-PrOH = 72/28, detector: 254 nm, flow rate: 0.80 mL/min, 30 °C), $t_1 = 15.7$ min (major), $t_2 = 17.2$ min; HRMS (ESI) *m*/*z* Calculated for $C_{16}H_{16}N_2O_3[M+H]^+$ 285.1234, found 285.1235.

(+)-4-Cyclohexyl-3,4-dihydroquinazolin-2(1*H*)-one (2n): 42 mg, 91% yield, white solid, mp: 130-131 °C, new compound, 96% ee, $[\alpha]^{20}_{D} = +14.3$ (*c* 0.40, MeOH), $R_f = 0.33$ (dichloromethane

/methanol = 15/1); ¹H NMR (400 MHz, DMSO-d₆) δ 8.97 (s, 1H), 7.12–7.08 (m, 1H), 7.03–6.95 (m, 2H), 6.87–6.83 (m, 1H), 6.77–6.75 (m, 1H), 4.10 (t, *J* = 3.8 Hz, 1H), 1.66 (t, *J* = 9.0 Hz, 2H), 1.61–1.48 (m, 3H), 1.44–1.36 (m, 1H), 1.19–0.90 (m, 5H); ¹³C NMR (100 MHz, DMSO-d₆) δ 154.8, 138.6, 127.9, 127.2,

121.0, 120.8, 113.8, 58.6, 46.8, 28.6, 27.1, 26.4, 26.2, 26.1; Enantiomeric excess was determined by HPLC (IC column, *n*-Hexane/*i*-PrOH = 92/08, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 21.1 \text{ min (major)}, t_2 = 22.5 \text{ min; HRMS (ESI)} m/z$ Calculated for $C_{14}H_{19}N_2O [M+H]^+$ 231.1492, found 231.1492.

(+)-4-Isopropyl-3,4-dihydroquinazolin-2(1*H*)-one (20): 35 mg, 92% yield, white solid, the known compound,¹⁰ 86% ee, $[\alpha]^{20}_{D}$ = +20.2 (*c* 0.56, MeOH), R_f = 0.40 (dichloromethane/

methanol = 15/1); ¹H NMR (400 MHz, CDCl₃) δ 8.09 (s, 1H), 7.17–7.13 (m, 1H), 7.03 (d, J = 7.4 Hz, 1H), 6.95 (t, J = 7.4 Hz, 1H), 6.73 (d, J = 7.8 Hz, 1H), 5.61 (s, 1H), 4.36 (t, J = 3.4 Hz, 1H), 2.00–1.91 (m, 1H), 0.99 (d, J = 6.8 Hz, 3H), 0.87 (d, J = 6.8 Hz, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 155.5, 136.7, 128.1, 126.6, 121.9,

120.3, 114.2, 60.0, 36.7, 18.5, 16.0; Enantiomeric excess was determined by HPLC (IC column, *n*-Hexane/*i*-PrOH = 90/10, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), $t_1 = 13.2$ min (major), $t_2 = 14.4$ min.

4. Asymmetric Hydrogenation at Gram Scale



A mixture of $[Ir(cod)Cl]_2$ (15.1 mg, 0.0225 mmol) and (*R*)-SegPhos (30.2 mg, 0.0495 mmol) in tetrahydrofuran (3.0 mL) was stirred at r.t. for 15 min in a glovebox, then BCDMH (108.7 mg, 0.45 mmol) and substrate **1a** (1.000 g, 4.5 mmol) together with tetrahydrofuran (13 mL) were added and the mixture was stirred for a further 10 min. The hydrogenation was performed at 25 °C under hydrogen (600 psi) in a stainless steel autoclave for 36 h. After carefully releasing the hydrogen, saturated aqueous sodium bicarbonate (10 mL) was added into the mixture and stirred for 10-15 min. The mixture was extracted with dichloromethane three times and the combined organic extract was dried over anhydrous sodium sulfate. After filtration, the filtrate was concentrated in *vacuo* and further purification was performed by a silica gel column with dichloromethane/methanol as eluent to give the desired product (*S*)-**2a** 0.931 g in 92% yield and 97% ee.

5. Synthesis of Bioactive Moelucules



The above synthetic methodology has been used as key step for facile syntheses of bioactive molecules. For example, (*S*)-**2a** could be converted into chiral thiourea (*S*)-**3a** with P_2S_5 in 67% yield, which is the Eg5 inhibitor.¹³ (*S*)-SDZ 267-489 **3b**, a serum HDL cholesterol raising agent,¹⁴ could be also synthesized in two steps from chiral (*S*)-**2a**.



The Synthesis of Eg5 Inhibitor: A mixture of (S)-2a (44.9 mg, 0.2 mmol, >99 ee) and P_2S_5 (38.3 mg, 0.2 mmol) in *p*-xylene (3 mL) was heated at 140 °C for 11 h under a nitrogen atmosphere. After being cooled to room temperature, the mixture was concentrated in *vacuo* and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give the desired product (S)-3a.

(*S*)-4-Phenyl-3,4-dihydroquinazoline-2(*1H*)-thione (3a): 32 mg, 67% yield, white solid, the known compound, $^{13,14} > 99\%$ ee, $[\alpha]^{20}{}_{D} = -160.41$ (*c* 0.24, MeOH), $R_f = 0.55$ (hexanes/ethyl acetate = 3/1); ¹H NMR (400 MHz, CDCl₃) δ 9.30 (s, 1H), 7.39-7.28 (m, 6H), 7.21-7.14 (m, 1H), 6.99-6.90 (m, 2H), 6.82 (d, J = 7.6 Hz, 1H), 5.66 (s, 1H); ¹³C NMR (100 MHz, CDCl₃) δ 174.6, 141.7, 133.5, 129.1, 128.8, 128.7, 127.5, 127.2, 124.3, 120.8, 114.4, 59.0; Enantiomeric excess

was determined by HPLC (OD-H column, *n*-Hexane/*i*-PrOH = 80/20, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), t₁ = 7.5 min (major), t₂ = 9.6 min.



The Synthesis of (S)-SDZ 267-489: A mixture of (S)-**3a** (48.1 mg, 0.20 mmol) and sodium bromide (4.1 mg, 0.04 mmol) was suspended in *i*-propanol (3 mL) and heated to 75 °C under a nitrogen atmosphere. To the rapidly stirred mixture was added chloroacetone (21 uL, 0.26 mmol) and the reaction was stirred at 75 °C for 3 h. After being cooled to room temperature, the mixture was concentrated in *vacuo* and further purification was performed by a silica gel column eluted with hexanes/ethyl acetate to give the desired product **3b** (S)-SDZ 267-489.

(*S*)-3-Methyl-5-phenyl-5*H*-thiazolo[2,3-*b*]quinazoline (3b): 46 mg, 83% yield, pale yellow solid, the known compound,¹⁴ 99% ee, $[\alpha]^{20}{}_{D} = -117.85$ (*c* 0.70, MeOH), $[lit.^{14} [\alpha]^{25}{}_{D} = -181.1$ (*c* 1.0, MeOH)], $R_f = 0.30$ (hexanes/ethyl acetate = 1/1); ¹H NMR (400 MHz, CDCl₃) δ 7.30-7.21 (m, 5H), 7.16-7.09 (m, 1H), 7.08-7.03 (m, 1H), 7.00-6.95 (m, 1H), 6.92-6.85 (m, 1H), 6.18 (s, 1H), 5.69 (s, 1H), 1.93 (s, 3H); ¹³C NMR (100 MHz, CDCl₃) δ 163.5, 143.6, 141.0, 134.6, 129.3, 128.8, 128.3, 126.6, 125.2, 123.4, 123.2, 121.3, 96.9, 61.0, 14.3; Enantiomeric excess was determined by HPLC (IC column, *n*-Hexane/*i*-PrOH = 80/20, detector: 254 nm, flow rate: 1.0 mL/min, 30 °C), t₁ = 19.1 min, t₂ = 20.5 min (major).

6. The Determination of Absolute Configuration

4-Phenyl-3,4-dihydroquinazolin-2(1*H*)-one (-)-**2a** was recrystallized in dichloromethane and *n*-hexane, optically pure product (> 99% ee) could be obtained. Then, a crystal was grown from dichloromethane and diethyl ether, which is suitable for X-ray diffraction analysis. The structure in **Figure S1** shows that the absolute configuration of (-)-**2a** is (4*S*). [CCDC 1480606] contains the structure and supplementary crystallographic data. These data can be obtained free of charge from the Cambridge Crystallographic Data Centre *via* www. ccdc.com.ac.uk/data_request/cif.



Figure S1. X-ray Crystallographic Analysis of (4S)-(-)-2a

7. References

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8. Copy of NMR and HPLC

















-22.14





13C NMR GF-4-46B in DMSO-d6







































13C NMR GF-4-49C IN DMSO-d6



~55.50





~56.61 ~55.55

13C NMR GF-4-45C IN DMSO-d6



¹³C NMR (100 MHz, DMSO-d₆)







13C NMR GF-4-49D in DMSO-d6









13C NMR GF-4-45D IN DMSO-d6



--56.41















































S53

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ009726.D Sample Name: GF-4-24E

Acq. Operator	:	1	
Acq. Instrument	:	Instrument 1 Location : Vial 1	
Injection Date	:	1/17/2016 9:00:50 AM	
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M	
Last changed	:	1/17/2016 6:46:33 AM by j	
		(modified after loading)	
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M	
Last changed	:	8/15/2018 8:57:38 PM	
		(modified after loading)	
Sample Info	:	OD-H, Hexane/iPrOH = 90/10, 1.0 mL/min, 30 oC, 254 nm	



Area Percent Report _____



Instrument 1 8/15/2018 8:57:48 PM



Page 1 of 1

*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZN000674.D Sample Name: GF-4-49E

Acq. Operator	:					
Acq. Instrument	:	Instrument 1	Location		Vial	1
Injection Date	:	3/20/2016 10:35:00 AM				
Acq. Method	:	C:\CHEM32\1\METHODS\DEF_LC.M				
Last changed	:	3/20/2016 10:12:51 AM				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	8/4/2018 2:26:00 PM				
		(modified after loading)				
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/m	nin, 30 o	с,	254 1	m



	Å1	cea Percent	t Report		
Sorted By Multiplier: Dilution: Use Multiplier « I	: Dilution H	Signal : : Factor with	1.0000 1.0000 h ISTDs		
<pre>>ignal 1: VWD1 A, Peak RetTime Type # [min]</pre>	Wavelengt Width [min] m	Area Area MAU *s	Height [mAU]	Area %	о НŅ ŅН
Signai 1: VWDI A, Peak RetTime Type # [min] 1 16.872 VB 2 20.712 BB	Wavelengt Vidth [min] m - 0.5676 5 0.6995	Area Area mAU *s 5212.28223 65.46461	Height [mAU] 141.17508 1.21299	Area % 98.7596 1.2404	

Instrument 1 8/4/2018 2:26:06 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010027.D Sample Name: GF-4-44A+-

Acq. Operator	:	†			
Acq. Instrument	:	Instrument 1	Location	:	Vial l
Injection Date	:	3/13/2016 8:41:15 AM			
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M			
Last changed	:	3/13/2016 6:08:33 AM by j			
		(modified after loading)			
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M			
Last changed	:	8/4/2018 1:36:54 PM			
		(modified after loading)			
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/m	nin, 30 o	С,	254 nm



Area Percent Report

Sorted By Multiplier: Dilution: Sample Amount: Use Multiplier & D	: Signal : : : vilution Factor w:	1.0000 1.0000 1.00000 [ng/ul] th ISTDs	(not used in calc.)
Signal 1: VWD1 A, Peak RetTime Type # [min] 1 12.828 EV 2 14.646 VBA	Wavelength=254 nm Width Area [min] mAU *s 0.4345 3381.4973 0.6183 3385.5810	Height Are [mAU] % 	a HN NH
Totals :	6767.0783	37 199.63944	(+/-)-2b

*** End of Report ***

Instrument 1 8/4/2018 1:37:04 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZ010048.D Sample Name: GF-4-44A

Acq. Operator	:	i				
Acq. Instrument	:	Instrument 1	Location	n :	Vial .	1
Injection Date	:	3/15/2016 10:59:31 AM				
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M				
Last changed	:	3/15/2016 8:50:00 AM by j				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M				
Last changed	:	8/4/2018 1:51:29 PM				
		(modified after loading)				
Sample Info	:	OD-H, H/1-PrOH = 90/10, 1.0 mL/m	nin, 30 (ьC,	254 m	m



 Area Percent Report

 Sorted By
 :
 Signal

 Multiplier:
 :
 1.0000

 Sample Amount:
 :
 1.0000

 Signal Amount:
 :
 1.0000

 Signal I:
 VWD1 A, Wavelength=254 nm
 (not used in calc.)

 Peak RetTime Type Width
 Area
 Height
 Area

 # [min]
 [min] mAU
 *s
 [mAU]
 *

 1
 12.6006 BB
 0.4194 1842.30322
 68.36278 96.0135
 I

 2
 14.711 BB
 0.4734 76.49199
 2.26079
 3.9865

 Totals :
 1918.79521 70.62357
 (-)-2b

Instrument 1 8/4/2018 1:51:36 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010019.D Sample Name: GF-4-44B+-

Acq. Operator	:	i		
Acq. Instrument	:	Instrument l Location : Vial 1		
Injection Date	:	3/13/2016 6:09:37 AM		
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M		
Last changed	:	3/13/2016 6:08:33 AM by j		
		(modified after loading)		
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M		
Last changed	:	8/4/2018 1:41:33 PM		
		(modified after loading)		
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 254 nm		



Area Percent Report

Sorted By	:	Sign	al		
Multiplier:		:	1.0000		
Dilution:		:	1.0000		
Sample Amount:		:	1.00000	[ng/ul]	(not used in calc.)
Use Multiplier	& Dilution	Factor	with ISTDs		



(+/-)-2c

*** End of Report ***

Instrument 1 8/4/2018 1:41:36 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\VZ010023.D Sample Name: GF-4-44B

	=						
Acq. Operator	÷	i					
Acq. Instrument	:	Instrument 1	Locatio	n	:	Vial	1
Injection Date	:	3/13/2016 7:22:35 AM					
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed	:	3/13/2016 6:08:33 AM by j					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	8/4/2018 1:53:47 PM					
		(modified after loading)					
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/	min, 30	оC	,	254 1	m



-----Area Percent Report _____ Sorted By Signal : : 1.0000 : 1.0000 : 1.00000 [ng/ul] (not used in calc.) Multiplier: Dilution: Sample Amount: Use Multiplier & Dilution Factor with ISTDs Signal 1: VWD1 A, Wavelength=254 nm 0 HN NH



*** End of Report ***

Instrument 1 8/4/2018 1:53:52 PM

Page 1 of 1

(-)-2c

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010031.D Sample Name: GF-4-44C+-

Acq. Operator	:	1		
Acq. Instrument	:	Instrument l Location : Vial 1		
Injection Date	:	3/13/2016 1:03:03 PM		
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M		
Last changed	:	3/13/2016 1:00:26 PM by j		
		(modified after loading)		
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M		
Last changed	:	8/4/2018 1:45:12 PM		
		(modified after loading)		
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 254 nm		



HN

'NH

(+/-)-2d

Page 1 of 1

Area Percent Report

Sorted By	:	Sig	nal		
Multiplier:		:	1.0000		
Dilution:		:	1.0000		
Sample Amount:		:	1.00000	[ng/ul]	(not used in calc.)
Use Multiplier &	Dilution	Factor	with ISTDs		



Instrument 1 8/4/2018 1:45:19 PM

*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZO10033.D Sample Name: GF-4-44C

:	i						
:	Instrument l L	ocation	. :	Vial 1			
:	3/13/2016 1:37:57 PM						
:	C:\HPCHEM\1\METHODS\DEF LC.M						
:	3/13/2016 1:00:26 PM by j						
	(modified after loading)						
:	C:\CHEM32\1\METHODS\DEF_LC11.M						
:	8/4/2018 1:48:34 PM						
	(modified after loading)						
:	OD-H, H/i-PrOH = 90/10, 1.0 mL/mi	n, 30 c	с,	254 nm			
		: 1 : 1 Instrument 1 : 3/13/2016 1:37:57 PM : C:\HPCHEWN\1METHODS\DEF LC.M : 3/13/2016 1:00:36 PM by 1 (modified after loading) : C:\CHEM33\1METHODS\DEF_LC11.M : 8/4/2018 1:48:34 PM (modified after loading) : OD-H, H/1-PrOH = 90/10, 1.0 mL/mi	: 1 : 1 Location : 3/13/2016 1:37:57 PM : C:\HPCHEN\1\METHODS\DEF LC.M : 3/13/2016 1:00:26 PM by j (modified after loading) : C:\CHENS2\1\METHODS\DEF_LC11.M : 8/4/2018 1:48:34 PM (modified after loading) : 0D-H, H/1-PrOH = 90/10, 1.0 mL/min, 30 c	: 1 : 1 Location : : 3/13/2016 1:37:57 PM : C:\HPCHEN\1\METHODS\DEF LC.M : 3/13/2016 1:00:26 PM hb j (modified after loading) : C:\CHEMS2\1\METHODS\DEF_LC11.M : 8/4/2018 1:48:34 PM (modified after loading) : 00-H, H/:1-FtOH = 90/10, 1.0 mL/min, 30 oC,	<pre>: 1 : 1 : Instrument 1</pre>		



-----Area Percent Report ______ Sorted By Signal : : 1.0000 : 1.0000 : 1.00000 [ng/ul] (not used in calc.) Multiplier: Dilution: Sample Amount: Use Multiplier & Dilution Factor with ISTDs Signal 1: VWD1 A, Wavelength=254 nm HN NH 1 18.513 BB 0.8599 1.48586e4 262.49072 97.4400 2 22.550 BB 1.0069 390.37512 5.00846 2.5600 Totals : 1.52490e4 267.49918

(-)-2d

*** End of Report ***

Instrument 1 8/4/2018 1:48:38 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZN000617.D Sample Name: GF-4-48B(+/-)

Acq. Operator	:							
Acq. Instrument	:	Instrument l Location : Vial 1						
Injection Date	:	3/17/2016 4:33:01 PM						
Acq. Method	:	C:\CHEM32\1\METHODS\DEF_LC.M						
Last changed	:	3/17/2016 4:28:44 PM						
		(modified after loading)						
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M						
Last changed	:	8/4/2018 2:14:46 PM						
		(modified after loading)						
Sample Info	:	OD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 254nm						





Data File C:\CHEM32\1\DATA\ZHOU-16\GF-2222\YZN000618.D Sample Name: GF-4-48B Acq. Instrument : Instrument 1 Location : Vial 1 Injection Date : 3/17/2016 6:11:39 PM Acq. Method : C:\CHEM32\1\METHODS\DFF_LC.M Last changed : 3/17/2016 6:04:47 PM (modified after loading) Analysis Method : C:\CHEM32\1\METHODS\DFF_LL.M Last changed : 8/4/2018 2:13:43 PM

(modified after loading) Sample Info : OD-H, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 254mm



-----Area Percent Report _ Signal Sorted By . Multiplier: : 1.0000 : 1.0000 Dilution: Use Multiplier & Dilution Factor with ISTDs \cap Signal 1: VWD1 A, Wavelength=254 nm HN NH Peak RetTime Type Width Area Height Area # [min] [min] mAU *s [mAU] % 1 13.126 VB 0.4811 1.53130e4 486.03174 97.8136 2 16.794 BB 0.7309 342.29477 7.01452 2.1864 Totals : 1.56553e4 493.04626 (-)-2e _____ *** End of Report ***

Instrument 1 8/4/2018 2:14:50 PM

Page 1 of 1

Instrument 1 8/4/2018 2:13:49 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010091.D Sample Name: GF-4-49C(+/-)

Acq. Operator	:	i						
Acq. Instrument	:	Instrument 1	Location	:	Vial 1			
Injection Date	:	3/19/2016 11:46:55 AM						
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M						
Last changed	:	3/19/2016 11:30:07 AM by j						
		(modified after loading)						
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M						
Last changed	:	8/4/2018 2:18:45 PM						
		(modified after loading)						
Sample Info	:	IC, H/i-PrOH = 80/20, 0.8 mL/mi	n, 30 oC,	2	54 nm			



Àrea Percent Report

Sorted By Multiplier: Dilution: Sample Amount: Use Multiplier	: & Dilution F	Sigmal : : : actor wit	1.0000 1.0000 1.00000 [ng h ISTDs	/ul] (not	t used in calc.)	
Signal 1: VWD1 Peak RetTime Ty # [min] 	A, Wavelengt pe Width [min] m 	h=254 nm Area AU *s 881.99121 875.23975	Height [mAU] 254.20068 147.92250	Area * 50.0287 49.9713	HN	O NH
Totals :	1	.17572e4	402.12318		`	OM∈ (+/-)-2f
	*	** End of	Report ***			

Data File C:\CHEM32\l\DATA\ZHOU-16\GF-2222\Y2010092.D Sample Name: GF-4-49C Acq. Operator : 1 Acq. Instrument : Instrument 1 Intection Date : 3/19/2016 12:28:38 PM Acq. Method : C:\FFPCHEM\L\METHODS\DEF LC.M Last changed : 3/19/2016 11:30:07 AM by 1 (modified after Loading) Analysis Method : C:\FHEN3L\METHODS\DEF_LC1L.M Last changed : 8/4/2018 2:17:25 PM







Instrument 1 8/4/2018 2:18:49 PM

Page 1 of 1

Instrument 1 8/4/2018 2:17:28 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010056.D
Sample Name: GF-4-45C(+/-)

Acq. Operator	:	1						
Acq. Instrument	:	Instrument 1 Location : Vial 1						
Injection Date	:	3/16/2016 3:01:16 AM						
Acq. Method	:	C:\HPCHEN\1\METHODS\DEF LC.M						
Last changed	:	3/16/2016 1:50:49 AM by j						
		(modified after loading)						
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M						
Last changed	:	8/4/2018 1:58:19 PM						
		(modified after loading)						
Sample Info	:	OD-3, H/i-PrOH = 90/10, 1.0 mL/min, 30 oC, 254 nm						



-----Area Percent Report -----Sorted Bv Sional

Sorted By	:	Sign	nal		
Multiplier:		:	1.0000		
Dilution:		:	1.0000		
Sample Amount:		:	1.00000	[ng/ul]	(not used in calc.)
Use Multiplier @	Dilution	Factor	with ISTDs		

Signal 1: VWD1 A,	Wavelen	gth=254 nm		
Peak RetTime Type # [min]	Width [min]	Area mAU *s	Height [mAU]	Area %
1 23.441 BV	0.8212	3888.69189	71.62736	49.0323
2 25.982 VB	0.7489	4042.18286	83.82679	50.9677
Totals :		7930.87476	155.45415	



*** End of Report ***

Instrument 1 8/4/2018 1:58:24 PM

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010055.D Sample Name: GF-4-45C

	=					
Acq. Operator	÷	i				
Acq. Instrument	:	Instrument 1	Location	:	Vial	1
Injection Date	:	3/16/2016 2:28:08 AM				
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M				
Last changed	:	3/16/2016 1:50:49 AM by j				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	8/4/2018 1:57:12 PM				
		(modified after loading)				
Sample Info	:	OD-3, H/i-PrOH = 90/10, 1.0 mL/m	in, 30 o	С,	254	nm



-----Area Percent Report _____ Sorted By Signal . Multiplier: 1.0000 : Dilution: . 1.0000 Sample Amount: : 1.00000 [ng/ul] (not used in calc.) Use Multiplier & Dilution Factor with ISTDs Signal 1: VWD1 A, Wavelength=254 nm \cap Peak RetTime Type Width Area Height Area

#	[min]		[min]	mAU *	s	[mAU]	*	
1	23.981	BV	0.7460	462.76	947	9.1	10235	4.4282	
2	25.732	VB	0.8208	9987.68	652	181.1	17450	95.5718	
Total	ls :			1.04505	e4	190.2	27685		



*** End of Report ***

Instrument 1 8/4/2018 1:57:17 PM

Page 1 of 1

(-)-2g

`OMe

HN

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010098.D Sample Name: GF-4-49D(+/-)

Acq. Operator	:	1						
Acq. Instrument	:	Instrument 1	Location	:	Vial l			
Injection Date	:	3/20/2016 2:05:49 AM						
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M						
Last changed	:	3/20/2016 1:55:38 AM by j						
		(modified after loading)						
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M						
Last changed	:	8/4/2018 2:23:00 PM						
		(modified after loading)						
Sample Info	:	IA, H/i-PrOH = 75/25, 0.9 mL/min	, 30 oC,	2	54 nm			



Area Percent Report _____ Signal Sorted By . : 1.0000 : 1.0000 : 1.0000 [ng/ul] (not used in calc.) Multiplier: Dilution: Sample Amount: Use Multiplier & Dilution Factor with ISTDs 0 Signal 1: VWD1 A, Wavelength=254 nm HN NH 1 11.858 BB 0.3148 2285.31616 107.60020 49.8263 2 20.745 BB 0.5658 2301.24805 61.74747 50.1737

4586.56421 169.34767

OMe (+/-)-2h

OMe

*** End of Report ***

Instrument 1 8/4/2018 2:23:06 PM

Totals :

Page 1 of 1

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZO10104.D Sample Name: GF-4-49D

	==			==:		
Acq. Operator	:	i				
Acq. Instrument	:	Instrument 1	Location	:	Vial	1
Injection Date	:	3/20/2016 7:37:40 AM				
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M				
Last changed	:	3/20/2016 7:17:27 AM by j				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	8/4/2018 2:21:35 PM				
		(modified after loading)				
Sample Info	:	IA, H/i-PrOH = 75/25, 0.9 mL/mi:	n, 30 oC,	2.	54 nm	
	Acq. Operator Acq. Instrument Injection Date Acq. Method Last changed Analysis Method Last changed Sample Info	Acq. Operator : Acq. Instrument : Intection Date : Acq. Method : Last changed : Last changed : Sample Info :	Acd. Operator : 1 Acd. Instrument : Instrument 1 Intection Date : 3/20/2016 7:37:40 AM Acq. Method : C:\HPCHEN\\\METHODS\DEF LC.M Last changed : 3/20/2016 7:17:27 AM by 1 (modified after loading) Analysis Method : C:\CHEM32\\METHODS\DEF_LCLI.M Last changed : 8/4/2018 2:21:35 PM (modified after loading) Sample Info : IA, H/i-Fr0H = 75/25, 0.9 mL/mi;	Acc. Operator : 1 Acc. Instrument : Instrument 1 Location Intection Date : 3/20/2015 7:37:40 AM Acc. Method : C:\HFCHEM\1\METHODS\DEF LC.M Last changed : 3/20/2016 7:17:27 AM by 1 (modified after loading) Analysis Method : C:\CHEM32\1\METHODS\DFF_LCll.M Last changed : 8/4/2018 2:21:35 PM (modified after loading) Sample Info : IA, Mi-FrOH = 75/25, 0.9 mL/min, 30 of,	Acd. Operator : 1 Acd. Instrument : Instrument 1 Intection Date : 3/20/2016 7:37:40 AM Acq. Method : C:\HPCHEM\1\METHODS\DEF LC.M Last changed : 3/20/2016 7:17:27 AM by j (modified after loading) Analysis Method : C:\CHEM32\1\METHODS\DEF_LC11.M Last changed : 8/4/2018 2:21:35 PM (modified after loading) Sample Info : IA, H/i-FPCM = 75/25, 0.9 mL/min, 30 oC, 2:	Acc. Operator : 1 Acc. Instrument : Instrument 1 Injection Date : 3/20/2015 7:37:40 AM Acq. Method : C:\HFCHEM\1\METHODS\DEF LC.M Last changed : 3/20/2016 7:17:27 AM by 1 (modified after loading) Analysis Method : C:\CHEM32\1\METHODS\DEF_LC11.M Last changed : 8/4/2018 2:21:35 PM (modified after loading) mathematical after loading) (modified after loading) (modified after loading) (modified after loading) (modified after loading)





Instrument 1 8/4/2018 2:21:39 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010061.D Sample Name: GF-4-45D(+/-)

Acq. Operator	:	†					
Acq. Instrument	:	Instrument l Location : Vial	1				
Injection Date	:	3/16/2016 7:12:01 AM					
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed	:	3/16/2016 7:08:44 AM by j					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	8/4/2018 2:02:18 PM					
		(modified after loading)					
Sample Info	:	OD-3, H/i-PrOH = 90/10, 0.9 mL/min, 30 oC, 254 m	m				



Area Percent Report

Sorted By	:	Signal			
Multiplier:		:	1.0000		
Dilution:		:	1.0000		
Sample Amount:		:	1.00000	[ng/ul]	(not used in calc.)
Use Multiplier	& Dilution	Factor wi	th ISTDs		



*** End of Report ***

 Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZO10062.D Sample Name: GF-4-45D

```
Acq. Operator : 1

Acq. Instrument : Instrument 1

Injection Date : 3/16/2016 7:39:23 AM

Acq. Method : C:\HFP(HEM)L\HETHODS\DEF LC.M

Last changed : 3/16/2016 7:08:44 AM by j

(modified after loading)

Analysis Method : C:\(HEM32)L\HETHODS\DEF LC1.M

Last changed : 8/4/2018 2:00:59 PM

(modified after loading)

Sample Info : 0D-3, H/i-FP(H = 90/10, 0.9 mL/min, 30 oC, 254 nm
```





*** End of Report ***

Instrument 1 8/4/2018 2:02:26 PM

Page 1 of 1

Instrument 1 8/4/2018 2:01:09 PM

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZ010060.D Sample Name: GF-4-45E(+/-)

Acq. Operator :	i						
Acq. Instrument :	Instrument 1	Location : Vial 1					
Injection Date :	3/16/2016 6:46:04 AM						
Acq. Method :	C:\HPCHEM\1\METHODS\DEF LC.M						
Last changed :	3/16/2016 5:04:34 AM by j						
	(modified after loading)						
Analysis Method :	C:\CHEM32\1\METHODS\DEF_LC11.M						
Last changed :	8/4/2018 2:05:49 PM						
	(modified after loading)						
Sample Info :	OD-3, H/i-PrOH = 90/10, 1.0 mL/	min, 30 oC, 254 nm					



-----Area Percent Report _____

Sorted By Multiplier: Dilution: Sample Amount: Use Multiplier &	: Dilution	Siqn : : Factor	nal 1.0000 1.0000 1.00000 with ISTDs	[ng/ul]	(not used in calc.)	
---	---------------	--------------------------	--	---------	---------------------	--

*** End of Report ***

Signal 1: VWD1 A,	Wavelength	=254 nm		
Peak RetTime Type	Width	Area	Height	Area
# [min]	[min] mA	U *s	[mAU]	%
1 14.969 BV	0.4300 1.	28553e4	439.52631	41.7410
2 16.626 VB	0.4412 1.	79425e4	603.96655	58.2590
Totals :	3.	07979e4	1043.49286	



Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010059.D Sample Name: GF-4-45E

Acq. Operator : j	i					
Acq. Instrument : I	Instrument 1	Location : Vial l				
Injection Date : 3	3/16/2016 6:03:00 AM					
Acq. Method : 0	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed : 3	3/16/2016 5:04:34 AM by j					
((modified after loading)					
Analysis Method : C	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed : 8	8/4/2018 2:04:37 PM					
	(modified after loading)					
Sample Info : 0	OD-3, H/i-PrOH = 90/10, 1.0 mL/m	in, 30 oC, 254 nm				



1	Area Percent Report		
			•
Sorted By :	Signal		
Multiplier:	: 1.0000		
Dilution:	: 1.0000		
Sample Amount:	: 1.00000 [no	/ull (not used in (alc.)
Use Multiplier & Dilution	Factor with ISTDs		•
·····			
Signal 1: VMD1 A. Naveleng	th=254 nm		0
bighti i. tobi ii, odocicii,	,		Ŷ
Peak DetTime Tyme Width	Area Height	Area	L L
# [min] [min]	mall *s [mall]	*	HN´ `NH
// [min] [min]		!	
1 15 419 377 0 3792	171 20261 6 67853	1 5167	\land
2 16 591 VP 0.4261	1 1116464 292 12121	00 4022	
2 10.301 VB 0.4301	1.1110464 303.12131	50.4033	
Tetels	1 12976-4 290 20094		
IUCAIS :	1.120/084 309./9904		~ ~ F
			(-)-2i
			()-)
	+++ Fud of Douout +++		
	···· End of Report ····		

Instrument 1 8/4/2018 2:05:52 PM

Page 1 of 1

Instrument 1 8/4/2018 2:04:43 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010276.D Sample Name: GF-4-67B-Rac

Acq. Operator	:	1					
Acq. Instrument	:	Instrument l Location : Vial 1					
Injection Date	:	4/4/2016 9:18:36 AM					
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed	:	4/4/2016 9:16:30 AM by j					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	8/4/2018 2:39:35 PM					
		(modified after loading)					
Sample Info	:	OD-H, H/i-PrOH = 95/05, 0.7 mL/min, 30 oC, 254 nm					



Area Percent Report

Sorted By : Multiplier: Dilution: Sample Amount: Use Multiplier & Dilutio	Signal : 1.0000 : 1.0000 : 1.00000 [ng/ul] (not used n Factor with ISTDs	in calc.)
Signal 1: VWD1 A, Wavele Peak RetTime Type Width # [min] [min]	ngth=254 nm Area Height Area mAU *s [mAU] %	
1 43.413 BB 1.268 2 47.895 BB 1.683	- 4 2325.61353 26.65312 51.8248 6 2161.83643 17.23040 48.1752	
Totals :	4487.44995 43.88352	(+/-)-2k
	*** End of Report ***	





Instrument 1 8/4/2018 2:39:40 PM

Page 1 of 1

Instrument 1 8/4/2018 2:42:13 PM

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZN000873.D Sample Name: GF-4-67C-Rac

Acq. Operator	:					
Acq. Instrument	:	Instrument 1	Location	:	Vial l	
Injection Date	:	4/3/2016 1:53:59 PM				
Acq. Method	:	C:\CHEM32\1\METHODS\DEF LC.M				
Last changed	:	4/3/2016 1:49:11 PM				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M				
Last changed	:	8/4/2018 2:45:25 PM				
		(modified after loading)				
Sample Info	:	OD-H, H/i-PrOH = 90/10, 0.7mL/m	in, 30oC,	2	54 nm	



Area Percent Report Sorted By Signal . Multiplier: : 1.0000 : 1.0000 Dilution: Use Multiplier & Dilution Factor with ISTDs 0 Signal 1: VWD1 A, Wavelength=254 nm HN NH Peak RetTime Type Width Area Height Area # [min] [min] mAU *s [mAU] ÷. 1 21.391 BV 0.6320 4353.11182 106.33298 49.9281 2 23.097 VB 0.9703 4365.64697 67.34595 50.0719 Totals : 8718.75879 173.67893 (+/-)-21 _____ *** End of Report ***

Data File C:\CHEM32\1\DATA\2H0U-16\GF-22222\YZN000874.D Sample Name: GF-4-67C Acr. Operator : Acq. Instrument : Instrument 1 Intection Date : 4/3/2016 2:25:22 PM Acq. Method : C:\CHEM32\1\METHODS\DFF_LC.M Last changed : 4/3/2016 2:22:49 PM (modified after loading) Analysis Method : C:\CHEM32\1\METHODS\DFF_LC1.M Last changed : G:\CHEM32\1\METHODS\DFF_LC1.M Last changed : G:\CHEM32\1\METHODS\DFF_LC1.M Last changed : G:\CHEM32\1\METHODS\DFF_LC1.M Sample Info : 0D-H, H/1-PrOH = 90/01,0.7mL/min, 300C, 254 nm





Instrument 1 8/4/2018 2:45:34 PM

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Instrument 1 8/4/2018 2:50:04 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010265.D Sample Name: GF-4-67A rac

Acq. Operator	:	i				
Acq. Instrument	:	Instrument 1	Location	:	Via.	11
Injection Date	:	4/4/2016 3:47:59 AM				
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M				
Last changed	:	4/4/2016 3:46:29 AM by j				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	8/4/2018 2:36:06 PM				
		(modified after loading)				
Sample Info	:	IC, H/i-PrOH = 72/28, 0.8 mL/mi	n, 30 oC,	2	54 m	m



Area Percent Report

- 1	Sorted By	:	Signal				
1	Multiplier:		:	1.0000			
1	Dilution:		:	1.0000			
-	Sample Amount:			1.00000 [nd	r/ull (not)	used in calc.)	
1	Use Multinlier & D	ilution	Factor with	n TSTDs	,	,	
							0
							Ŷ
	Firmel L. WHD L A	Horrolone	sth-254 mm				Ц
	Jighai I. Vobi A,	wavereni	jui-234 im				HN [^] NH
	Deels DetTine Trme	Ui d+h		Unisht	Awaa		
	reak Reclime Type	forder 1	ALea	Teronic J	Area		
	# [min]	נשדעו	IIAU "S	[IIIAO]			
	1 15.761 VV	0.4175	2272.89722	84.40331	49.9074		Ľ // Ľ //
	2 17.125 VB	0.4707	2281.33325	74.51803	50.0926	MeO	Ý Ý
	Fotals :		4554.23047	158.92133			OMe
							(./) 2m
							(+/-)-2111
			*** End of	Report ***			
				-			





Instrument 1 8/4/2018 2:36:14 PM

Page 1 of 1

Instrument 1 8/4/2018 2:37:22 PM

Data File C:\CHEM32\1\DATA\ZH0U-16\GF-22222\YZ010110.D Sample Name: GF-4-52C(+/-)

Acq. Operator	:	i					
Acq. Instrument	:	Instrument 1	Locat	ion	: 1	/ial	1
Injection Date	:	3/21/2016 12:35:48 PM					
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed	:	3/21/2016 12:34:33 PM by j					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	8/4/2018 2:32:53 PM					
		(modified after loading)					
Sample Info	:	IC, H/i-PrOH = 92/08, 1.0 mL/mi	n, 30	oC,	254	1 nm	



Area Percent Report

Sorted By : Multiplier: Dilution: Sample Amount: Use Multiplier & Dilution Signal 1: VWD1 A, Wavele:	Signal : 1.0000 : 1.0000 : 1.00000 [ng/ul] (not used in cald n Factor with ISTDs ngth=254 nm	
Peak RetTime Type Width # [min] [min] 	Area Height Area mAU *s [mAU] % 4 2074.29028 63.81608 49.5694 6 2110.33154 59.59352 50.4306 4184.62183 123.40960 123.40960	(+/-)-2n

*** End of Report ***

Data File C:\CHEH32\l\DATA\ZHOU-16\GF-22222\Y2010112.D Sample Name: GF-4-52C Acq. Operator : 1 Acq. Instrument : Instrument 1 Location : Vial 1 Infection Date : 3/22/2016 1:16:37 AM Acq. Method : C:\HPCHEM\L\METHODS\DEF LC.M Last changed : 3/22/2016 1:13:42 AM by j (modified after 1 oading)

Analysis Method : C:\CHEM32\1\METHODS\DEF_LC11.M Last changed : 8/4/2018 2:34:19 PM (modified after loading) Sample Info : IC, H/-PrOH = 92/08, 1.0 mL/min, 30 oC, 254 nm



Area Percent Report _ Signal Sorted By : Multiplier: 1.0000 : 1.0000 Dilution: -Sample Amount: . 1.00000 [ng/ul] (not used in calc.) Use Multiplier & Dilution Factor with ISTDs 0 Signal 1: VWD1 A, Wavelength=254 nm HN NH 1 21.067 BV 0.4934 7212.34521 227.62305 97.8503 2 22.507 VB 0.5368 158.44650 4.55595 2.1497 (+)-2n Totals : 7370.79172 232.17900

*** End of Report ***

Instrument 1 8/4/2018 2:32:58 PM

Page 1 of 1

Instrument 1 8/4/2018 2:34:23 PM

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZ010108.D Sample Name: GF-4-52B(+/-)

Acq. Operator	:	i				
Acq. Instrument	:	Instrument 1	Location	:	Vial l	
Injection Date	:	3/21/2016 11:50:24 AM				
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M				
Last changed	:	3/21/2016 11:33:47 AM by j				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M				
Last changed	:	8/4/2018 2:28:21 PM				
		(modified after loading)				
Sample Info	:	IC, H/i-PrOH = 90/10, 1.0 mL/mi	n, 30 oC,	2	54 nm	



Area Percent Report

Sorted By Multiplier: Dilution: Sample Amount: Use Multiplier &	: Dilution F	Signal : : actor with	1.0000 1.0000 1.00000 [n 1.STDs	g/ul]	(not used in	calc.) O
Signal 1: VWD1 A	, Wavelengt	h=254 nm				HN
Peak RetTime Typ # [min] 1 13.441 BV 2 14.616 VB	e Width [min] m - - 0.2979 4 0.3254 4	Area AU *s .160.84570 .174.30029	Height [mAU] 218.48232 199.66394	Area % 49.919 50.080	 93 17	
Totals :	8	335.14600	418.14626			(+/-)-20
						=

*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-16\GF-22222\YZO10111.D Sample Name: GF-4-52B

Acq. Operator	:	i					
Acq. Instrument	:	Instrument 1	Locat	ion	:	Vial	1
Injection Date	:	3/22/2016 12:52:04 AM					
Acq. Method	:	C:\HPCHEM\1\METHODS\DEF LC.M					
Last changed	:	3/22/2016 12:09:58 AM by j					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	8/4/2018 2:31:03 PM					
		(modified after loading)					
Sample Info	:	IC, H/i-PrOH = 90/10, 1.0 mL/min	1, 30	oC,	25	54 nm	



	Ar	ea Percen	t Report				
Sorted By	:	Signal					
Multiplier:		:	1.0000				
Dilution:		:	1.0000				
Sample Amount:		:	1.00000	[ng/ul]	(not used	in calc.)	
Use Multiplier 🤅	Dilution F	'actor wit	h ISTDs				•
Gi	Horrelengt	h-254 mm					Ļ
Peak RetTime Type	, wavelengt e Width	Area	Height	Area			
Peak RetTime Type # [min]	, waverengc e Width [min] m	Area AU *s	Height [mAU]	Area %	1		
Peak RetTime Type # [min] 1 13.213 BB	, wavelengt e Width [min] m 	Area AU *s 	Height [mAU] 222.339	Area % 22 93.07	1		HN' NH
Peak RetTime Type # [min] 1 13.213 BB 2 14.375 BB	wavelengt e Width [min] m 	Area AU *s 070.75049 302.89621	Height [mAU] 222.339 14.903	Area * 22 93.07 15 6.92	 45 55		HN' NH

*** End of Report ***

Instrument 1 8/4/2018 2:28:27 PM

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Instrument 1 8/4/2018 2:31:10 PM

Data File C:\CHEM32\1\DATA\ZHOU-18\YZNO10811.D Sample Name: zz-4-78(+-)

Acq. Operator	:					
Acq. Instrument	:	Instrument 1	Location	:	-	
Injection Date	:	11/20/2018 10:31:17 PM				
Acq. Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	11/20/2018 10:24:12 PM				
		(modified after loading)				
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M				
Last changed	:	11/29/2018 10:32:35 AM				
		(modified after loading)				
Sample Info	:	OD-H, Hexane/i-PrOH =80/20, 1.0	mL/min, 3	30 ol	C, 254 ni	m



Sorted	By		:	Signal			
Multip	lier:			:	1.0000		
Diluti	on:			:	1.0000		
Use Mu	ltiplie	er « D	ilution	Factor wit	h ISTDs		
Signal	1: VWD)1 A,	Waveleng	rth=254 nm			
Signal	1: VWI)1 A,	Wavelenq	gth=254 nm			
Signal Peak R	l: VWD etTime)1 A, Type	Wavelen Width	yth=254 nm Area	Height	Area	
Signal Peak R #	l: VWI etTime [min])l A, Type	Waveleng Width [min]	yth=254 nm Area mAU *s	Height [mAU]	Area %	
Signal Peak R # -	l: VUI etTime [min])1 A, Type	Waveleng Width [min]	gth=254 nm Area mAU *s	Height [mAU]	Area %	
Signal Peak R # - 1	1: VWD etTime [min] 7.502)1 Å, Type VV	Waveleng Width [min] 0.2382	gth=254 nm Area mAU *s 1269.43982	Height [mAU] 	Area % 49.1545	
Signal Peak R # - 1 2	1: VWC etTime [min] 7.502 9.589)1 A, Type VV VB	Waveleng Width [min] 0.2382 0.3630	gth=254 nm Area mAU *s 1269.43982 1313.10974	Height [mAU] 80.97906 54.42619	Area * 49.1545 50.8455	
Signal Peak R # - 1 2	1: VWD [min] 7.502 9.589	Dl A, Type VV VB	Waveleng Width [min] 0.2382 0.3630	gth=254 nm Area mAU *s 1269.43982 1313.10974	Height mAU 80.97906 54.42619	Area * 49.1545 50.8455	

*** End of Report ***

Data File C:\CHEM32\1\DATA\ZHOU-18\YZNO10955.D Sample Name: zz-4-87-2

Acq. operator							
Acq. Instrument	:	Instrument 1	Location	:	-		
Injection Date	:	12/4/2018 4:56:43 PM					
Acq. Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M					
Last changed	:	12/4/2018 4:39:39 PM					
		(modified after loading)					
Analysis Method	:	C:\CHEM32\1\METHODS\DEF LC11.M					
Last changed	:	12/7/2018 6:32:56 PM					
		(modified after loading)					
Sample Info	:	OD-H, Hexane/i-PrOH = 80/20, 1.	0 mL/min,	30	οC,	254	nm







Instrument 1 11/29/2018 10:32:43 AM

Page 1 of 1

(+/-)-3a

HN

Instrument 1 12/7/2018 6:32:58 PM

Data File G:\广收液相\SIG1001993.D Sample Name: zz-4-83(+-)

	-	
Acq. Operator	:	
Acq. Instrument	:	仪器 1 Location : Vial 91
Injection Date	:	11/27/2018 10:52:36 AM
		Inj Volume : 5.000 µl
Acq. Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M
Last changed	:	11/27/2018 10:45:12 AM
		(modified after loading)
Analysis Method	:	C:\CHEM32\1\METHODS\DEF_LC11.M
Last changed	:	12/7/2018 6:38:18 PM
		(modified after loading)
Sample Info	:	IC, n-hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 254 nm



Area Percent Report

Sorted By : Signal Multiphier: : 1.0000 Dilution: : 1.0000 Use Multiphier & Dilution Factor with ISTDs

Signal 1: VWDl A, Wavelength=254 nm Peak RetTime Type Width Area Height Area # [min] [min] [mAU³5] [mAU] ²

1	18.763	VV	0.4192	2429.51733	88.81035	49.5783
2	20.381	VB	0.4684	2470.84814	80.88632	50.4217
Total	s :			4900.36548	169.69667	

*** End of Report ***

Instrument 1 12/7/2018 6:38:24 PM

Page 1 of 1

Ρh

(+/-)-3b

Me

Data File G:\广收液相\SIG1002049.D Sample Name: zz-3-88

Acq. Operator :	
Acq. Instrument :	仪器 1 Location : Vial 91
Injection Date :	12/5/2018 7:13:25 PM
	Inj Volume : 5.000 µl
Acq. Method :	C:\CHEM32\1\METHODS\DEF_LC11.M
Last changed :	12/5/2018 6:45:39 PM
	(modified after loading)
Analysis Method :	C:\CHEN32\1\METHODS\DEF LC11.M
Last changed :	12/7/2018 6:37:42 PM
	(modified after loading)
Sample Info :	IC, n-hexane/i-PrOH = 80/20, 1.0 mL/min, 30 oC, 254 nm



Area Percent Report

Sorted By	:	Signal	
Multiplier:		:	1.0000
Dilution:		:	1.0000
Use Multiplier	& Dilution	Factor with	i ISTDs

Signal 1: VWD1 A, Wavelength=254 nm

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.074	VV	0.4447	9.80341	3.43762e-1	0.4773
2	20.463	VB	0.4807	2044.18359	65.38577	99.5227
Total	s :			2053.98700	65.72953	



*** End of Report ***

Instrument 1 12/7/2018 6:37:46 PM